# Intensity Frontier Common Offline Documentation: art Workbook and Users Guide

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#### art Glossary

abstraction the process by which data and programs are defined with a representa-

tion similar in form to its meaning (semantics), while hiding away the implementation details. A system can have several abstraction layers whereby different meanings and amounts of detail are exposed to the programmer (adapted from Wikipedia's entry for "Abstraction (com-

puter science)".

analyzer module an art module that may read information from the current event but that

may not add information to it; e.g., a module to fill histograms or make

printed output

API Application Programming Interface

art The art framework (art is not an acronym) is the software framework

developed for common use by the Intensity Frontier experiments to

develop their offline code and non-real-time online code

art module see *module* 

art path a FHiCL sequence of *art* moduleLabels that specifies the work the job

will do

artdaq a toolkit that lives on top of *art* for building high-performance event-

building and event-filtering systems; this toolkit is designed to support efficient use of multi-core computers and GPUs. A technical paper on

artdag can be found at.

bash a UNIX shell scripting language that is used by some of the support

scripts in the workbook exercises

boost a class library with new functionality that is being prototyped for in-

clusion in future C++ standards

build system turns source code into object files, puts them into a shared library, links

them with other libraries, and may also run tests, deploy code to pro-

duction systems and create some documentation.

buildtool a Fermilab-developed tool (part of **cetbuildtools**) to compile, link and

run tests on the source code of the Workbook

catch See *exception* in a C++ reference

cetbuildtools a build system developed at Fermilab

CETLIB a utility library used by *art* (developed and maintained by the *art* team)

to hold information that does not fit naturally into other libraries

class The C++ programming language allows programmers to define program-

specific data types through the use of *classes*. Classes define types of data structures and the functions that operate on those data structures. Instances of these data types are known as *objects*. Other object ori-

ented languages have similar concepts.

CLHEP a set of utility classes; the name is an acronym for a Class Library for

**HEP** 

collection

configuration see run-time configuration

const member function a member function of a class that does not change the value of

non-mutable data members: see mutable data member

constructor a function that (a) shares an identifier with its associated class, and (b)

initializes the members of an object instantiated from this class

DAQ data aquisition system

data handling

Data Model see Event Data Model

data product Experiment-defined class that can represent detector signals, recon-

structed data, simulated events, etc. In *art*, a *data product* is the smallest unit of information that can be added to or retrieved from an event.

data type See *type* 

declaration (of a class) the portion of a class that specifies its type, its name, and any data

members and/or member functions it has

destructor a function that (a) has the same identifier as its associated class but

prefaced with a tilde ( $\sim$ ), and (b) is used to deallocate memory and do other cleanup for a class object and its class members when the object

is destroyed

Doxygen a system of producing reference documentation based on comments in

source code

ED a prefix used in *art* (e.g., for module types) meaning *event-data* 

EDAnalyzer see analyzer module

EDFilter see filter module

EDOutput see *output module* 

EDProducer see producer module

EDSource see source module

Event In HEP there are two notions of the word *event* that are in common use;

see event (unit of information) or event (interaction). In this documen-

tation suite, unless otherwise indicated, we mean the former.

Event (interaction) An event (unit of data) may contain more than one fundamental in-

teraction; the science goal is always to identify individual fundamental interactions and determine their properties. It is common to use the word *event* to refer to one of the individual fundamental interactions. In the near detector of a high-intensity neutrino experiment, for example, there may be multiple neutrino interactions within the unit of time that defines a single *event (unit of information)*. Similarly, in a colliding-beam experiment, an *event (unit of information)* corresponds

to the information from one beam crossing, during which time there may be multiple collisions between beam particles.

Event (unit of information) In the general HEP sense, an *event* is a set of raw data associated in time, plus any information computed from the raw data; *event* may also refer to a simulated version of same. Within *art*, the representation of an *event* (*unit of information*) is the classs art::Event, which is the smallest unit of information that *art* can process. An art::Event contains an event identifier plus an arbitrary number of *data-products*; the information within the *data-products* is intrinsically experiment dependent and is defined by each experiment. For bookkeeping convenience, *art* groups events into a heirarchy: a *run* contains zero or more *subRuns* and a *subRun* contains zero or more events.

Event Data Model (EDM) Representation of the data that an experiment collects, all the derived information, and historical records necessary for reproduction of result

within an *art* job, the set of steps to perform in order to execute the per-event functions for each event that is read in, including steps for begin/end-job, begin/end-run and begin/end-subRun

all of the data products in an experiment's files; plus the metadata that accompanies them. The HEP software community has adopted the word *event-data* to refer to the software details of dealing with the information found in *events*, whether the events come from experimental data or simulations.

event-data file a collective noun to describe both data files and files of simulated events

exception, to throw a mechanism in C++ (and other programming languages) to stop the current execution of a program and transfer control up the call chain; also called *catch* 

experiment code see user code

event loop

event-data

external product for a given experiment, this is a software product that the experiment's software (within the *art* framework) does not build, but that it uses; e.g., ROOT, Geant4, etc. At Fermilab external products are managed

by the in-house UPS/UPD system, and are often called *UPS products* or simply *products*.

FermiGrid a batch system for submitting jobs that require large amounts of CPU

time

FHICL Fermilab Hierarchical Configuration Language (pronounced "fickle"),

a language developed and maintained by the *art* team at Fermilab to support run-time configuration for several projects, including *art* 

FHiCL-CPP the C++ toolkit used to read FHiCL documents within art

filter module an *art* module that may alter the flow of processing modules within an

event; it may add information to the event

framework (art) The art framework is an application used to build physics programs

by loading physics algorithms, provided as plug-in modules; each experiment or user group may write and manage its own modules. *art* also provides infrastructure for common tasks, such as reading input, writing output, provenance tracking, database access and run-time con-

figuration.

framework (generic) an abstraction in which software providing generic functionality can

be selectively changed by additional user-written code, thus providing application-specific software (significantly abbreviated from Wikipedia's entry for "software framework"); note that the actual functionality provided by any given framework, e.g., *art*, will be tailored to the given

needs.

free function a function without data members; it knows only about agruments passed

to it at run time; see function and member function

Geant4 a toolkit for the simulation of the passage of particles through matter,

developed at CERN. http://geant4.cern.ch/

git a source code management system used to manage files in the art

Workbook; similar in concept to the older CVS and SVN, but with

enhanced functionality

handle a type of smart pointer that permits the viewing of information inside

a data product but does not allow modification of that information; see *pointer*, data product

IF Intensity Frontier

ifdh sam a UPS product that allows *art* to use *SAM* as an external run-time agent

that can deliver remote files to local disk space and can copy output files to tape. The first part of the name is an acronym for Intensity Frontier

Data Handling.

implementation the portion of C++ code that specifies the functionality of a declared

data type; where as a struct or class declaration (of a data type) usually resides in a *header* file (.h or .hh), the implementation usually resides in a separate source code file (.cc) that "#includes" the header file

instance see *instantiation* 

instantiation the creation of an object instance of a class in an OOP language; an

instantiated object is given a name and created in memory or on disk

using the structure described within its class declaration.

jobsub-tools a UPS product that supplies tools for submitting jobs to the Fermigrid

batch system and monitoring them.

Kerberos a single sign-on, strong authentication system required by Fermilab for

access to its computing resources

kinit a command for obtaining Kerberos credentials that allow access to Fer-

milab computing resources; see Kerberos

member function (also called *method*) a function that is defined within (is a *member* of)

a class; they define the behavior to be exhibited by instances of the associated class at program run time. At run time, member functions have access to data stored in the instance of the class with they are associated, and are thereby able to control or provide access to the state

of the instance.

message facility a UPS product used by art and experiments' code that provides facili-

ties for merging messages with a variety of severity levels, e.g., infor-

mational, error, and so on; see also *mf* 

message service

method see *member function* 

mf a namespace that holds classes and functions that make up the message

facility used by art and by experiments that use art; see message facility

module a C++ class that obeys certain rules established by art and whose source

code file gets compiled into a shared object library that can be dynamically loaded by *art*. An *art* module "plugs into" a processing stream and performs a specific task on units of data obtained using the Event Data Model, independent of other running modules. See also *module*-

Label

module\_type

moduleLabel a user-defined identifier whose value is a parameter set that *art* will use

to configure a module; see module and parameter set

Monte Carlo method a class of computational algorithms that rely on repeated random

sampling to obtain numerical results; i.e., by running simulations many times over in order to calculate those same probabilities heuristically just like actually playing and recording your results in a real casino

situation: hence the name (Wikipedia)

mutable data member The

namespace a container within a file system for a set of identifiers (names); usu-

ally grouped by functionality, they are used to keep different subsets of code distinguishable from one another; identical names defined within different namespaces are disambiguated via their namespace prefix

ntuple an ordered list of n elements used to describe objects such as vectors or

tables

object an instantiation of any data type, built-in types (e.g., int, double, float)

or class types; i.e., a location range in memory containing an instantia-

tion

object-oriented language a programming language that supports OOP; this usually means

support for classes, including public and private data and functions

object-oriented programming (OOP) a programming language model organized around

*objects* rather than procedures, where *objects* are quantities of interest that can be manipulated. (In contrast, programs have been viewed historically as logical procedures that read in data, process the data and produce output.) Objects are defined by *classes* that contain attributes (data fields that describe the objects) and associated procedures. See *C++ class*; *object*.

OOP see object oriented programming

output module an art module that writes data products to output file(s); it may select

a subset of data products in a subset of events; an art module contains

zero or more output modules

parameter set a C++ class, defined by FHICL-CPP, that is used to hold run-time

configuration for *art* itself or for modules and services instantiated by *art*. In a FHiCL file, a parameter set is represented by a FHiCL *table*;

see *table* 

path a generic word based on the UNIX concept of PATH that refers to a

colon-separated list of directories used by art when searching for vari-

ous files (e.g., data input, configuration, and so on)

physics in *art*, *physics* is the label for a portion of the run-time configuration

of a job; this portion contains up to five sections, each labeled with a

reserved

pointer a variable whose value is the address of (i.e., that points to) a piece

of information in memory. A native C++ pointer is often referred to as a *bare pointer*. *art* defines different sorts of *smart pointers* (or *safe pointers*) for use in different circumstances. One commonly used type

of smart pointer is called a handle.

process\_name a parameter to which the user assigns a mnemonic value identifying

the physics content of the associated FHiCL parameter set (i.e., the parameters used in the same FHiCL file). The process\_name value is

embedded into every data product created via the FHiCL file.

producer module an art module that may read information from the current event and

may add information to it

product See either *external* product or *data* product

redmine an open source, web-based project management and bug-tracking tool

used as a repository for art code and related code and documentation

ROOT an HEP data management and data presentation package used by art

and supported by CERN; *art* is designed to allow output of event-data to files in ROOT format, in fact currently it is the only output format

that art implements

ROOT files There are two types of ROOT files managed by art: (1) event-data

output files, and (2) the file managed by TFileService that holds user-

defined histograms, ntuples, trees, etc.

run a period of data collection, defined by the experiment (usually delin-

eates a period of time during which certain running conditions remain

unchanged); a run contains zero or more subRuns

run-time configuration (processing-related) structured documents describing all process-

ing aspects of a single job including the specification of parameters and

workflow; in art it is supplied by a FHiCL file; see FHiCL

safe pointer see *pointer* 

SAM (Sequential data Access via Metadata) a Fermilab-supplied product

that provides the functions of a file catalog, a replica manager and some

functions of a batch-oriented workflow manager

scope

sequence (in FHiCL) one or more comma-separated FHiCL values delimited by square

brackets (

• • •

) in a FHiCL file is called a *sequence* (as distinct from a *table*)

service in art, a singleton-like object (type) whose lifetime and configuration

are managed by *art*, and which can by accessed by module code and by

other services by requesting a service handle to that particular service.

The service *type* is used to provide geometrical information, conditions and management of the random number state; it is also used to implement some internal functionality. See also *T File Service* 

shared library

signature (of a function) the unique identifier of a C++ a function, which includes: (a)

its name, including any class name or namespace components, (b) the number and type of its arguments, (c) whether it is a member function, (d) whether it is a const function (Note that the signature of a function

does not include its return type.)

site As used in the *art* documentation, a *site* is a unique combination of

experiment and institution; used to refer to a set of computing resources configured for use by a particular experiment at a particular institution. This means that, for example, the Workbook environment on a Mu2e-owned computer at Fermilab will be different than that on an Mu2e-owned computer at LBL. Also, the Workbook environment on a Mu2e-owned computer at Fermilab will be different from that on an LBNE-

owned computer at Fermilab.

smart pointer see *pointer* 

source (refers to a *data* source) the name of the parameter set inside an FHiCL

file describing the first step in the workflow for processing an event; it reads in each event sequentially from a data file or creates an empty

event; see also source code; see also EDsource

source code code written in C++ (the programming language used with art) that

requires compilation and linking to create an executable program

source module an *art* module that can initiate an *art* path by reading in event(s) from

a data file or by creating an empty event; it is the first step of the pro-

cessing chain

standard library, C++ the C++ standard library of routines

std identifier for the namespace used by the C++ standard library

struct identical to a C++ class except all members are *public* (instead of *pri*-

vate) by default

subRun

a period of data collection within a run, defined by the experiment (it may delineate a period of time during which certain run parameters remain unchanged); a SubRun is contained within a run; a subRun contains zero or more events

table (in FHiCL)

a group of FHiCL definitions delimited by braces ({ ... }) is called a table; within art, a FHiCL table gets turned into an object called a parameter set. Consequently, a FHiCL table is typically called a parameter set. See parameter set.

template (C++)

Templates are a feature of C++ that allows for meta-programming. In practical terms, the coder can write an algorithm that is independent of type, as long as the type supports the features required by the algorithm. For example, there is a standard library "sort" algorithm that will work for any type that provides a way to determine if one object of the type is "less than" another object of the type.

**TFileService** 

an art service used by all experiments to give each module a ROOT subdirectory in which to place its own histograms, TTrees, and so on; see TTrees and ROOT

truth information One use of simulated events is to develop, debug and characterize the algorithms used in reconstruction and analysis. To assist in these tasks, the simulation code often creates data products that contain detailed information about the right answers at intermediate stages of reconstruction and analysis; they also write data products that allow the physicist to ask "is this a case in which there is an irreducible background or should I be able to do better?" This information is called the truth information, the Monte Carlo truth or the God's block.

**TTrees** 

a ROOT implementation of a tree; see *tree* and *ROOT* 

type

Variables and objects in C++ must be classified into types, e.g., builtin types (integer, boolean, float, character, etc.), more complex userdefined classes/structures and typedefs; see *class*, *struct*, and *typedef*. The word type in the context of C++ and art is the same as data type unless otherwise stated.

typedef A typedef is a different name, or an alias, by which a type can be identi-

fied. Type aliases can be used to reduce the length of long or confusing type names, but they are most useful as tools to abstract programs from

the underlying types they use (cplusplus.com).

UPS/UPD a Fermliab-developed system for distributing software products

user code experiment-specific and/or analysis-specific C++ code that uses the art

framework; this includes any personal code you write that uses art.

variable a storage location and an associated symbolic name (an identifier) which

contains some known or unknown quantity or information, a value. The variable name is the usual way to reference the stored value; this separation of name and content allows the name to be used independently

of the exact information it represents.

# Part I

## Introduction

## 1 How to Read this Documentation

The *art* document suite, which is currently in an alpha release form, consists of an introductory section and the first few exercises of the Workbook\*, plus a glossary and an index. There are also some preliminary (incomplete and unreviewed) portions of the Users Guide included in the compilation.

The Workbook exercises require you to download some code to edit, execute and evaluate. Both the documentation and the code it references are expected to undergo continual development throughout 2013 and 2014. The latest is always available at the *art* Documentation website. Chapter 11 tells you how to keep up-to-date with improvements and additions to the Workbook code and documentation.



## 1.1 If you are new to HEP Software...

Read Parts I and II (the introductory material and the Workbook) from start to finish. The Workbook is aimed at an audience who is familiar with (although not necessarily expert in) Unix, C++ and Fermilab's UPS product management system, and who understands the basic *art* framework concepts. The introductory chapters prepare the "just starting out" reader in all these areas.

## 1.2 If you are an HEP Software expert...

Read chapters 1, 2 and 3: this is where key terms and concepts used throughout the *art* document suite get defined. Skip the rest of the introductory material and jump straight

<sup>\*</sup>The Workbook is expected to contain roughly 35 exercises when complete.

into running Exercise 1 in Chapter 9 of the Workbook. Take the approach of: Don't need it? Don't read it.

## 1.3 If you are somewhere in between...

Read chapters 1, 2 and 3 and skim the remaining introductory material in Part I to glean what you need. Along with the experts, you can take the approach of: Don't need it? Don't read it.

## 2 Conventions Used in this Documentation

Most of the material in this introduction and in the Workbook is written so that it can be understood by those new to HEP computing; if it is not, please let us know (see Section 3.4)!

The first instance of each term that is defined in the glossary is written in *italics* followed by a  $\gamma$  (Greek letter gamma), e.g.,  $framework(\gamma)$ .

Unix commands that you must type are shown in the format unix command. Portions of the command for which you must substitute values are surrounded by angle brackets (< ... >, e.g., you would type your actual username when you see <username>).

When an example Unix command line would overflow the page width, this documentation will use a trailing backslash to indicate that the command is continued on the next line. For example:

```
mkdir -p \ ART\_WORKBOOK\_WORKING\_BASE/< username > / workbook-tutorial / \\ directory 1 / directory 2 / directory 3
```

This means that you should type the entire command on a single line if it fits, without typing the backslash, *or* on two lines *with* the backslash as the final character of the first line.

Step-by-step procedures that the reader is asked to follow are denoted in this way. Commands inside procedures are denoted as mkdir-p <mydir>.

Computer output from a command is shown as:

Part

#### command output

In some places it will be necessary for a paragraph or two to be written for experts. Such paragraphs will be marked with a "dangerous bends" symbol in the margin, as shown at right. Less experienced users can skip these sections on first reading and come back to them at a later time.



Occasionally, text will be called out to make sure that you don't miss it. Important or tricky terms and concepts will be marked with an "pointing finger" symbol in the margin, as shown at right.



Items that are even trickier will be marked with a "bomb" symbol in the margin, as shown at right. You really want to avoid the problems they describe.



Text that refers in particular to Fermilab-specific information is marked with a Fermilab picture, as shown at right.



Text that refers in particular to information about using *art* at non-Fermilab sites is marked with a "generic site" picture, as shown at right. A *site* is defined as a unique combination of experiment and institution, and is used to refer to a set of computing resources configured for use by a particular experiment at a particular institution.



Experiment-specific information will be kept to an absolute minimum; wherever it appears, it will be marked with an experiment-specific icon, e.g., the Mu2e icon at right.



# 3 Introduction to the art Event Processing Framework

#### 3.1 What is art and Who Uses it?

 $art(\gamma)$  is an event-processing  $framework(\gamma)$  developed and supported by the Fermilab Scientific Computing Division (SCD). The art framework is used to build physics programs by loading physics algorithms, provided as plug-in modules. Each experiment or user group may write and manage its own modules. art also provides infrastructure for common tasks, such as reading input, writing output, provenance tracking, database access and run-time configuration.

The initial clients of *art* are the Fermilab Intensity Frontier experiments but nothing prevents other experiments from using it as well. The name *art* is always written in *italic lower case*; it is not an acronym.

art is written in C++ and is intended to be used with user code written in C++. (*User code* includes experiment-specific code and any other user-written, non-art, non-external-product( $\gamma$ ) code.)

art has been designed for use in most places that a typical HEP experiment might require a software framework, including:

- high-level software triggers
- o online data monitoring
- o calibration
- o reconstruction

- o analysis
- o simulation

art is not designed for use in real-time environments, such as the direct interface with data-collection hardware.

The Fermilab SCD has also developed a related product named  $artdaq(\gamma)$ , a layer that lives on top of art and provides features to support the construction of data-acquisition  $(DAQ(\gamma))$  systems based on commodity servers. Further discussion of artdaq is outside the scope of this documentation; for more information consult the artdaq redmine site:

https://cdcvs.fnal.gov/redmine/projects/artdaq/wiki.

The design of *art* has been informed by the lessons learned by the many High Energy Physics (HEP) experiments that have developed C++ based frameworks over the past 20 years. In particular, it was originally forked from the framework for the CMS experiment, *cmsrun*.

Experiments using art are listed at the art Documentation website under "Experiments using art."

## 3.2 Why *art*?

In all previous experiments at Fermilab, and in most previous experiments elsewhere, infrastructure software (i.e., the framework, broadly construed – mostly forms of bookkeeping) has been written in-house by each experiment, and each implementation has been tightly coupled to that experiment's code. This tight coupling has made it difficult to share the framework among experiments, resulting in both great duplication of effort and mixed quality.

art was created as a way to share a single framework across many experiments. In particular, the design of art draws a clear boundary between the framework and the user code; the art framework (and other aspects of the infrastructure) is developed and maintained by software engineers who are specialists in the field of HEP infrastructure software; this provides a robust, professionaly maintained foundation upon which physicists can develop the code for their experiments. Experiments use art as an external package. Despite some

constraints that this separation imposes, it has improved the overall quality of the framework and reduced the duplicated effort.

## 3.3 C++ and C++11

In 2011, the International Standards Committee voted to approve a new standard for C++, called C++ 11.

Much of the existing user code was written prior to the adoption of the C++ 11 standard and has not yet been updated. As you work on your experiment, you are likely to encounter both code written the new way and code written the old way. Therefore, the Workbook will often illustrate both practices.

A very useful compilation of what is new in C++ 11 can be found at



https://cdcvs.fnal.gov/redmine/projects/gm2public/wiki/CPP2011

This reference material is written for advanced C++ users.

## 3.4 Getting Help

Please send your questions and comments to *art-users@fnal.gov*. More support information is listed at https://sharepoint.fnal.gov/project/ArtDoc-Pub/SitePages/Support.aspx.

## 3.5 Overview of the Documentation Suite

When complete, this documentation suite will contain several principal components, or *volumes*: the introduction that you are reading now, a Workbook, a Users Guide, a Reference Manual, a Technical Reference and a Glossary. At the time of writing, drafts exist for the Introduction, the Workbook, the Users Guide and the Glossary. The components in the documentation suite are illustrated in Figure 3.1.

Part

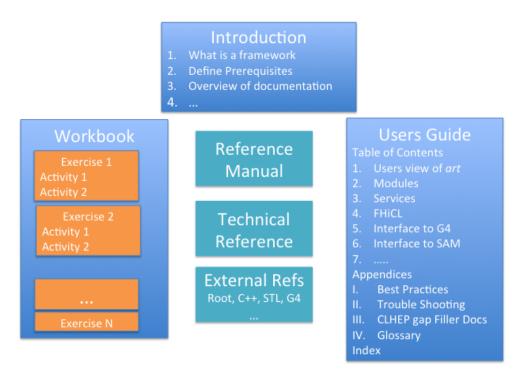


Figure 3.1: The principal components of the art documentation suite

#### 3.5.1 The Introduction

This introductory volume is intended to set the stage for using *art*. It introduces *art*, provides background material, describes some of the software tools on which *art* depends, describes its interaction with related software and identifies prerequisites for successfully completing the Workbook exercises.

#### 3.5.2 The Workbook

The Workbook is a series of standalone, self-paced exercises that will introduce the building blocks of the *art* framework and the concepts around which it is built, show practical applications of this framework, and provide references to other portions of the documentation suite as needed. It is targeted towards physicists who are new users of *art*, with the understanding that such users will frequently be new to the field of computing for HEP and to C++.

One of the Workbook's primary functions is training readers how and where to find more extensive documentation on both *art* and external software tools; they will need this information as they move on to develop and use the scientific software for their experiment.

The Workbook assumes some basic computing skills and some basic familiarity with the C++ computing language; Chapter 6 provides a tutorial/refresher for readers who need to improve their C++ skills.

The Workbook is written using recommended best practices that have become current since the adoption of C++ 11 (see Section 3.8).

Because *art* is being used by many experiments, the Workbook exercises are designed around a *toy* experiment that is greatly simplified compared to any actual detector, but it incorporates enough richness to illustrate most of the features of *art*. The goal is to enable the physicists who work through the exercises to translate the lessons learned there into the environment of their own experiments.

#### 3.5.3 Users Guide

The Users Guide is targeted at physicists who have reached an intermediate level of competence with art and its underlying tools. It contains detailed descriptions of the features of art, as seen by the physicists. The Users Guide will provide references to the  $external\ products(\gamma)$  on which art depends, information on how art uses these products, and as needed, documentation that is missing from the external products' own documentation.

#### 3.5.4 Reference Manual

The Reference Manual will be targeted at physicists who already understand the major ideas underlying *art* and who need a compact reference to the Application Programmer Interface ( $API(\gamma)$ ). The Reference Manual will likely be generated from annoted source files, possibly using  $Doxygen(\gamma)$ .

#### 3.5.5 Technical Reference

The Technical Reference will be targeted at the experts who develop and maintain *art*; few physicists will ever want or need to consult it. It will document the internals of *art* so that a broader group of people can participate in development and maintenance.

#### 3.5.6 Glossary

The glossary will evolve as the documentation set grows. At the time of writing, it includes definitions of *art*-specific terms as well as some HEP, Fermilab, C++ and other relevant computing-related terms used in the Workbook and the Users Guide.

## 3.6 Some Background Material

This section defines some language and some background material about the *art* framework that you will need to understand before starting the Workbook.

art Documentation

#### 3.6.1 Events and Event IDs

In almost all HEP experiments, the core idea underlying all bookkeeping is the  $event(\gamma)$ . In a triggered experiment, an event is defined as all of the information associated with a single trigger; in an untriggered, spill-oriented experiment, an event is defined as all of the information associated with a single spill of the beam from the accelerator. Another way of saying this is that an event contains all of the information associated with some time interval, but the precise definition of the time interval changes from one experiment to another \*. Typically these time intervals are a few nanoseconds to a few tens of mircoseconds. The information within an event includes both the raw data read from the Data Acquisition System (DAQ) and all information that is derived from that raw data by the reconstruction and analysis algorithms. An event is the smallest unit of data that art can process at one time.

In a typical HEP experiment, the trigger or DAQ system assigns an event identifier (event ID) to each event; this ID uniquely identifies each event, satisfying a critical requirement imposed by *art* that each event be uniquely identifiable by its event ID. This requirement also applies to simulated events.

The simplest event ID is a monotonically increasing integer. A more common practice is to define a multi-part ID and *art* has chosen to use a three-part ID, including:

- $\circ$  run( $\gamma$ ) number
- $\circ$  *subRun*( $\gamma$ ) number
- $\circ$  event( $\gamma$ ) number

There are two common methods of using this event ID scheme and *art* allows experiments to chose either:

1. When an experiment takes data, the event number is incremented every event. When some predefined condition occurs, the event number is reset to 1 and the subRun number is incremented, keeping the run number unchanged. This cycle repeats until some other predefined condition occurs, at which time the event number is reset to

<sup>\*</sup>There is a second, distinct, sense in which the word *event* is sometimes used; it is used as a synonym for a *fundamental interaction*; see the glossary entry for *event (fundamental interaction)*( $\gamma$ ). Within this documentation suite, unless otherwise indicated, the word *event* refers to the definition given in the main body of the text.

- 1, the subRun number is reset to 0 (0 not 1 for historical reasons) and the run number is incremented.
- 2. The second method is the same as the first except that the event number monontonically increases throughout a run and does not reset to 1 on subRun boundaries. The event number does reset to 1 at the start of each run.

art does not define what conditions cause these transitions; those decisions are left to each experiment. Typically experiments will choose to start new runs or new subRuns when one of the following happens: a preset number of events is acquired; a preset time interval expires; a disk file holding the ouptut reaches a preset size; or certain running conditions change.

art requires only that a subRun contain zero or more events and that a run contain zero or more subRuns.

When an experiment takes data, events read from the DAQ are typically written to disk files, with copies made on tape. The events in a single subRun may be spread over several files; conversely, a single file may contain many runs, each of which contains many subRuns.

#### 3.6.2 art Modules and the Event Loop

Users provide executable code to art in pieces called  $art \, modules(\gamma)^{\dagger}$  that are dynamically loaded as plugins and that operate on event data. The concept of reading events and, in response to each new event, calling the appropriate member functions of each module, is referred to as the  $event \, loop(\gamma)$ . The concepts of the  $art \, module$  and the  $event \, loop$  will be illustrated via the following discussion of how art processes a job.

The simplest command to run *art* looks like:

The argument to -c is the *run-time configuration file*( $\gamma$ ), a text file that tells one run of *art* what it should do. Run-time configuration files for *art* are written in the Fermilab

<sup>†</sup>Many programming languagues have an idea named *module*; the use of the term *module* by *art* and in this documentation set is an *art*-specific idea that will be developed through the first few chapters of the Workbook.

Hierarchical Configuration Language  $FHiCL(\gamma)$  (pronounced "fickle") and the filenames end in .fcl. As you progress through the Workbook, this language and the conventions used in the run-time configuration file will be explained; the full details are available in Chapter 25 of the Users Guide. (The run-time configuration file is often referred to as simply the *configuration file* or even more simply as just the *configuration*( $\gamma$ ).)

When *art* starts up, it reads the configuration file to learn what input files it should read, what user code it should run and what output files it should write. As mentioned above, an experiment's code (including any code written by individual experimenters) is provided in units called *art modules*. A module is simply a C++ class, provided by the experiment or user, that obeys a set of rules defined by *art* and whose *source*  $code(\gamma)$  file gets compiled into a shared  $object(\gamma)$  library that can be dynamically loaded by *art*. These rules will be explained as you work through the Workbook and they are summarized in Section 32.3



The code base of a typical experiment will contain many C++ classes. Only a small fraction of these will be modules; most of the rest will be ordinary C++ classes that are used within modules<sup>‡</sup>.

A user can tell *art* the order in which modules should be run by specifying that order in the configuration file. A user can also tell *art* to determine, on its own, the correct order in which to run modules; the latter option is referred to as *reconstruction on demand*.

Imagine the processing of each event as the assembly of a widget on an assembly line and imagine each module as a worker that needs to perform a set task on each widget. Each worker has a task that must be done on each widget that passes by; in addition some workers may need to do some start-up or close-down jobs. Following this metaphor, *art* requires that each module provide code that will be called once for every event and *art* allows any module to provide code that will be called at the following times:

- o at the start of the art job
- o at the end of the art job
- o at the start of each run
- o at the end of each run

 $<sup>\</sup>ddagger$  art defines a few other specialized roles for C++ classes; you will encounter these in Sections 3.6.4 and 3.6.5.

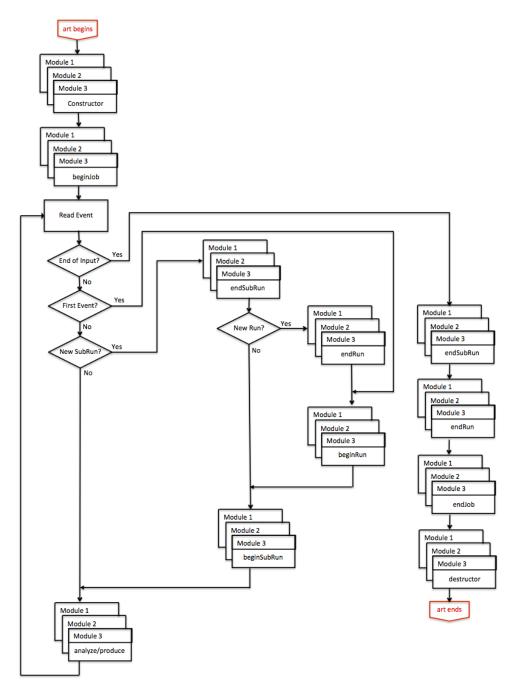
- o at the start of each SubRun
- o at the end of each SubRun

For those of you who are familiar with *inheritance* in C++, a module class (i.e., a "module") must inherit from one of a few different module *base classes*. Each module class must override one pure-virtual member function from the base class and it may override other virtual member functions from the base class.



After *art* completes its initialization phase (intentionally not detailed here), it executes the event loop. This is illustrated in Figure 3.2, which is described in the text below:

- 1. calls the  $constructor(\gamma)$  of every module in the configuration
- 2. calls the beginJob member function( $\gamma$ ) of every module that provides one
- 3. reads one event from the input source, and for that event
  - (a) determines if it is from a run different from that of the previous event (true for first event in loop)
  - (b) if so, calls the beginRun member function of each module that provides one
  - (c) determines if the event is from a subRun different from that of the previous event (true for first event in loop)
  - (d) if so, calls the beginSubRun member function of each module that provides one
  - (e) calls each module's (required) per-event member function
- 4. reads the next event and repeats the above per-event steps until it encounters a new subRun
- closes out the current subRun by calling the endSubRun member function of each module that provides one
- 6. repeats steps 4 and 5 until it encounters a new run
- 7. closes out the current run by calling the endRun member function of each module that provides one
- 8. repeats steps 3 through 7 until it reaches the end of the input source



**Figure 3.2:** Flowchart describing the *art* event loop for an input file that contains at least one event. *art* begins at the box in the upper left and ends at the box in the lower right. On the first event, the tests for new subRun and new run are true. Not all features of the event loop are shown, just those that you will encounter in the early parts of the *art* workbook. The case of a file with no events is not shown because it has many subcases and is not of general interest.

- 9. calls the endJob member function of each module that provides one
- 10. calls the  $destructor(\gamma)$  of each module

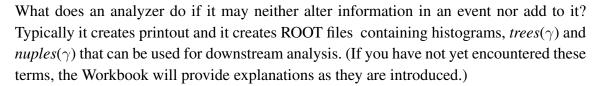
This entire set of steps comprises the event loop. One of *art*'s most visible jobs is controlling the event loop.

#### 3.6.3 Module Types

Every *art* module must be one of the following five types, which are defined by the ways in which they interact with each event and with the event loop:

- analyzer module( $\gamma$ ) May inspect information found in the event but may not add new information to the event; described in Chapter 28.
- **producer module**( $\gamma$ ) May inspect information found in the event and may add new information to the event; described in Chapter 27.
- *filter module*( $\gamma$ ) Same functions as a producer module but may also tell *art* to skip the processing of some, or all, modules for the current event; may also control which events are written to which output; described in Chapter 29.
- *source module*( $\gamma$ ) Reads events, one at a time, from some source; *art* requires that every *art* job contain exactly one source module. A source is often a disk file but other options exist and will be described in the Workbook and Users Guide.
- *output module*( $\gamma$ ) Reads an event from memory and writes it to an output; an *art* job may contain zero or more output modules. An ouptut is often a disk file but other options exist and will be described in the Workbook and in .

Note that no module may change information that is already present in an event.



Most novice users will only write analyzer modules and filter modules; readers with a little more experience may also write producer modules. The Workbook will provide examples of all three. Few people other than *art* experts and each experiment's software experts will



write source or output modules, however, the Workbook will teach you what you need to know about configuring source and output modules.

#### 3.6.4 art Data Products

This section introduces more ideas and terms dealing with event information that you will need as you progress through the Workbook.

The term  $data\ product(\gamma)$  is used in art to mean the unit of information that user code may add to an event or retrieve from an event. A typical experiment will have the following sorts of data products:

- 1. The DAQ system will package the raw data into data products, perhaps one or two data products for each major subsystem.
- 2. Each module in the reconstruction chain will create one or more data products.
- 3. Some modules in the analysis chain will produce data products; others may just make histograms and write information in non-*art* formats for analysis outside of *art*; they may, for example, write user-defined ROOT TTrees.
- 4. The simulation chain will usually create many data products. Some will be simulated event-data while others will describe the true properties of the simulated event. These data products can be used to study the response of the detector to simulated events; they can also be used to develop, debug and characterize the reconstruction algorithms.

Because these data products are intrinsically experiment-dependent, each experiment defines its own data products. In the Workbook, you will learn about a set of data products designed for use with the toy experiment. There are a small number of data products that are defined by *art* and that hold bookkeeping information; these will be described as you encounter them in the Workbook.



A data product is just a C++  $type(\gamma)$  (a class,  $struct(\gamma)$  or typedef) that obeys a set of rules defined by art; these rules are very different than the rules that must be followed for a class to be a module . A data product can be a single integer, an large complex class hierarchy, or anything in between.

Very often, a data product is a  $collection(\gamma)$  of some experiment-defined type. The C++

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standard libraries define many sorts of collection types; *art* supports many of these and also provides a custom collection type named cet::map\_vector. Workbook exercises will clarify the *data product* and *collection type* concepts.

#### 3.6.5 art Services

Previous sections of this Introduction have introduced the concept of C++ classes that have to obey a certain set of rules defined by art, in particular, modules in Section 3.6.2 and data products in Section 3.6.4.  $art services(\gamma)$  are yet another example of this.

In a typical *art* job, two sorts of information need to be shared among the modules. The first sort is stored in the data products themselves and is passed from module to module via the event. The second sort is not associated with each event, but rather is valid for some aggregation of events, subRuns or runs, or over some other time interval. Three examples of this second sort include the geometry specification, the conditions information and, for simulations, the table of particle properties.

To provide managed access to the second sort of information, *art* supports an idea named *art services* (again, shortened to *services*). Services may also be used to provide certain types of utility functions. Again, a service in *art* is just a C++ class that obeys a set of rules defined by *art*. The rules for services are different than those for modules or data products.

art implements a number of services that it uses for internal functions, a few of which you will encounter in the first couple of Workbook exercises. The message service( $\gamma$ ) is used by both art and experiment-specific code to limit printout of messages with a low severity level and to route messages to appropriate destinations. It can be configured to provide summary information at the end of the art job. The TFileService( $\gamma$ ) and the RandomNumberGenerator service are not used internally by art, but are used by most experiments. Experiments may also create and implement their own services.

After *art* completes its initialization phase and before it constructs any modules (see Section 3.6.2), it

<sup>§</sup>The phrase "conditions information" is the currently fashionable name for what was once called "calbration constants"; the name change came about because most calibration information is intrinsically time-dependent, which makes "constants" a poor choice of name.

- 1. reads the configuration to learn what services are requested
- 2. calls the constructor of each requested service

Once a service has been constructed, any code in any module can ask art for a smart  $pointer(\gamma)$  to that service and use the features provided by that service. Because services are constructed before modules, they are available for use by modules over the full life cycle of each module.

It is also legal for one service to request information from another service as long as the dependency chain does not have any loops. That is, if Service A uses Service B, then Service B may not use Service A, either directly or indirectly.



For those of you familiar with the C++ Singleton Design Pattern, an *art* service has some differences and some similarities to a Singleton. The most important difference is that the lifetime of a service is managed by *art*, which calls the constructors of all services at a well-defined time in a well-defined order. Contrast this with the behavior of Singletons, for which the order of initialization is undefined by the C++ standard and which is an accident of the implementation details of the loader. *art* also includes services under the umbrella of its powerful run-time configuration system; in the Singleton Design pattern this issue is simply not addressed.

#### 3.6.6 Shareable Libraries and art

When code is executed within the *art* framework, *art*, not the experiment, provides the main executable. The experiment provides its code to the *art* executable in the form of shareable object libraries that *art* loads dynamically at run time; these libraries are also called *dynamic load libraries* or *plugins* and their filenames are required to end in .so. For more information about shareable libraries, see Section 32.5.

#### 3.6.7 Build Systems and art

To make an experiment's code available to *art*, the source code must be compiled and linked (i.e., *built*) to produce shareable object libraries (Section 3.6.6). The tool that creates the .so files from the C++ source files is called a *build system*( $\gamma$ ).

Experiments that use art are free to choose their own build systems, as long as the system

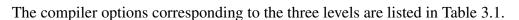
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**Table 3.1:** Compiler flags for the optimization levels defined by **cetbuildtools**; compiler options not related to optimization or debugging are not included in this table.

Name	flags
debug	-O0 -g
prof	-O3 -g -fno-omit-frame-pointer -DNDEBUG
opt	-O3 -DNDEBUG

follows the conventions that allow art to find the name of the .so file given the name of the module class. The Workbook will use a build system named cetbuildtools, which is a layer on top of  $cmake^{\P}$ .

The **cetbuildtools** system defines three standard compiler optimization levels, called "debug", "profile" and "optimized"; the last two are often abbreviated "prof" and "opt". When code is compiled with the "opt" option, it runs as quickly as possible but is difficult to debug. When code is compiled with the "debug" option, it is much easier to debug but it runs more slowly. When code is compiled with the "prof" option the speed is almost as fast as for an "opt" build and the most useful subset of the debugging information is retained. The "prof" build retains enough debugging information that one may use a profiling tool to identify in which functions the program spends most of its time; hence its name "profile". The "prof" build provides enough information to get a useful traceback from a core dump. Most experiments using *art* use the "prof" build for production and the "debug" build for development.





#### 3.6.8 External Products

As you progress through the Workbook, you will see that the exercises use some software packages that are part of neither *art* nor the toy experiment's code. The Workbook code, *art* and the software for your experiment all rely heavily on some external tools and, in order to be an effective user of *art*-based HEP software, you will need at least some familiarity with them; you may, in fact, need to become expert in some.

These packages and tools are referred to as *external products*( $\gamma$ ) (sometimes called simply *products*).

 $<sup>\</sup>P$ **cetbuildtools** is also used to build *art* itself.

An initial list of the external products you will need to become familiar with includes:

art the event processing framework

FHiCL the run-time configuration language used by art

CETLIB a utility library used by art

 $MF(\gamma)$  a message facility that is used by art and by (some) experiments that use art

ROOT an analysis, data presentation and data storage tool widely used in HEP

 $CLHEP(\gamma)$  a set of utility classes; the name is an acronym for Class Library for HEP

 $boost(\gamma)$  a class library with new functionality that is being prototyped for inclusion in future C++ standards

- gcc the GNU C++ compiler and run-time libraries; both the core language and the standard library are used by *art* and by your experiment's code.
- $git(\gamma)$  a source code management system that is used for the Workbook and by some experiments; similar in concept to the older CVS and SVN, but with enhanced functionality
- $cetbuildtools(\gamma)$  a Fermilab-developed external product that contains buildtool and related tools
- $UPS(\gamma)$  a Fermilab-developed system for accessing software products; it is an acronym for *Unix Product Support*.
- $UPD(\gamma)$  a Fermilab-developed system for distributing software products; it is an acronym for *Unix Product Distribution*.
- $jobusub\_tools(\gamma)$  tools for submitting jobs to the Fermigrid batch system and monitoring them.
- $ifdh\_sam(\gamma)$  allows art to use  $SAM(\gamma)$  as an external run-time agent that can deliver remote files to local disk space and can copy output files to tape. SAM is a Fermilab-supplied resource that provides the functions of a file catalog, a replica manager and some functions of a batch-oriented workflow manager

Any particular line of code in a Workbook exercise may use elements from, say, four or five of these packages. Knowing how to parse a line and identify which feature comes from

which package is a critical skill. The Workbook will provide a tour of the above packages so that you will recognize elements when they are used and you will learn where to find the necessary documentation.

For the *art* Workbook, external products are made available to your code via a mechanism called UPS, which will be described in Section 7. Many Fermilab experiments also use UPS to manage their external products; this is not required by *art* and you may choose to manage external products whichever way you prefer. UPS is, itself, just another external product. From the point of view of your experiment, *art* is an external product. From the point of view of the Workbook code, both *art* and the code for the toy experiment are external products.

Finally, it is important to recognize an overloaded word, *products*. When a line of documentation simply says *products*, it may be referring either to data products or to external products. If it is not clear from the context which is meant, please let us know (see Section 3.4).



#### 3.6.9 The Event-Data Model and Persistency

Section 3.6.4 introduced the idea of *art* data products. In a small experiment, a fully reconstructed event may contain on the order of ten data products; in a large experiment there may be hundreds.

While each experiment will define its own data product classes, there are many issues that are common to all data products in all experiments:

- 1. How does my module access data products that are already in the event?
- 2. How does my module publish a data product so that other modules can see it?
- 3. How is a data product represented in the memory of a running program?
- 4. How does an object in one data product refer to an object in another data product?
- 5. What metadata is there to describe each data product?

  Such metadata might include: which module created it; what was the run-time configuration of that module; what data products were read by that module; what was the code version of the module that created it?

6. How does my module access the metadata associated with a particular data product?

The answers to these questions form what is called the *Event-Data Model*( $\gamma$ ) (EDM) that is supported by the framework.

A question that is closely related to the EDM is: what technologies are supported to write data products from memory to a disk file and to read them from the disk file back into memory in a separate *art* job? A framework may support several such technologies. *art* currently supports only one disk file format, a ROOT-based format, but the *art* EDM has been designed so that it will be straightforward to support other disk file formats as it becomes useful to do so.

A few other related terms that you will encounter include:

- 1. transient representation: the in-memory representation of a data product
- 2. persistent representation: the on-disk representation of a data product
- 3. *persistency*: the technology to convert data products back and forth between their persistent and transient representations

#### 3.6.10 Event-Data Files

When you read data from an experiment and write the data to a disk file, that disk file is usually called a *data file*.

When you simulate an experiment and write a disk file that holds the information produced by the simulation, what should you call the file? The Particle Data Group has recommended that this not be called a "data file" or a "simulated data file;" they prefer that the word "data" be strictly reserved for information that comes from an actual experiment. They recommend that we refer to these files as "files of simulated events" or "files of Monte Carlo events" Note the use of "events," not "data."

This leaves us with a need for a collective noun to describe both data files and files of simulated events. The name in current use is *event-data files*( $\gamma$ ); yes this does contain the word "data" but the hyphenated word, "event-data", is unambiguous and this has become the standard name.

In HEP almost all simulations codes use *Monte Carlo*( $\gamma$ ) methods; therefore simulated events are often referred to as *Monte Carlo events* and the simulation process is referred to as *running the Monte Carlo*.

#### 3.6.11 Files on Tape

Many experiments do not have access to enough disk space to hold all of their event-data files, ROOT files and log files. The solution is to copy a subset of the disk files to tape and to read them back from tape as necessary.

At any given time, a snapshot of an experiment's files will show some on tape only, some on tape with copies on disk, and some on disk only. For any given file, there may be multiple copies on disk and those copies may be distributed across many  $sites(\gamma)$ , some at Fermilab and others at collaborating laboratories or universities.

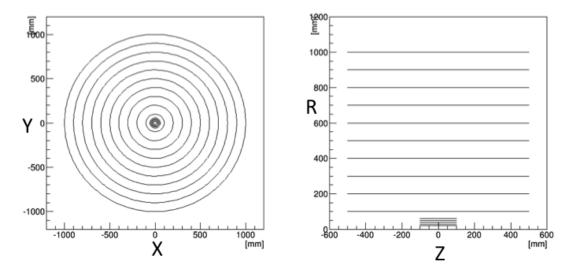
Conceptually, two pieces of software are used to keep track of which files are where, a *File Catalog* and a *Replica Manager*. One software package that fills both of these roles is called SAM, which is an acronym for "Sequential data Access via Metadata." SAM also provides some tools for Workflow management. SAM is in wide use at Fermilab and you can learn more about SAM at:

https://cdcvs.fnal.gov/redmine/projects/sam-main/wiki.

## 3.7 The Toy Experiment

The Workbook exercises are based around a made-up (toy) experiment. The code for the toy experiment is deployed as a UPS product named toyExperiment. The rest of this section will describe the physics content of toyExperiment; the discussion of the code in the toyExperiment UPS product will unfold in the Workbook, in parallel to the exposition of art.

The software for the toy experiment is designed around a toy detector, which is shown in Figure 3.3. The *toyExperiment* code contains many C++ classes: some modules, some data products, some services and some plain old C++ classes. About half of the modules are producers that individually perform either one step of the simulation process or one step of the reconstruction/analysis process. The other modules are analyzers that make histograms and ntuples of the information produced by the producers. There are also event display modules.



**Figure 3.3:** The geometry of the toy detector; the figures are described in the text. A uniform magnetic field of strength 1.5 T is oriented in the +z direction.

#### 3.7.1 Toy Detector Description

The toy detector is a central detector made up of 15 concentric shells, with their axes centered on the z axis; the left-hand part of Figure 3.3 shows an xy view of these shells and the right shows the radius vs z view. The inner five shells are closely spaced radially and are short in z; the ten outer shells are more widely spaced radially and are longer in z. The detector sits in a uniform magnetic field of 1.5 T oriented in the +z direction. The origin of the coordinate system is at the center of the detector. The detector is placed in a vacuum.

Each shell is a detector that measures  $(\varphi, z)$ , where  $\varphi$  is the azimuthal angle of a line from the origin to the measurement point. Each measurement has perfectly gaussian measurement errors and the detector always has perfect separation of hits that are near to each other. The geometry of each shell, its efficiency and resolution are all configurable at runtime.

All of the code in the toyExperiment product works in the set of units described in Table 3.2. Because the code in the Workbook is built on toyExperiment, it uses the same units. *art* itself is not unit-aware and places no constraints on which units your experiment may use.

Quantity Unit Length mm Energy MeV Time ns Plane Angle radian Solid Angle steradian Electric Charge Charge of the proton = +1Magnetic Field Tesla

**Table 3.2:** Units used in the Workbook

The first six units listed in Table 3.2 are the base units defined by the CLHEP SystemOfUnits package. These are also the units used by Geant4.



#### 3.7.2 Workflow for Running the Toy Experiment Code

The workflow of the toy experiment code includes five steps: three simulation steps, a reconstruction step and an analysis step:

- 1. event generation
- 2. detector simulation
- 3. hit-making
- 4. track reconstruction
- 5. analysis of the mass resolution

For each event, the event generator creates some signal particles and some background particles. The first signal particle is generated with the following properties:

- $\circ$  Its mass is the rest mass of the  $\phi$  meson; the event generator does not simulate a natural width for this particle.
- It is produced at the origin.
- $\circ$  It has a momentum that is chosen randomly from a distribution that is uniform between 0 and 2000 MeV/c.
- Its direction is chosen randomly on the unit sphere.

The event generator then decays this particle to  $K^+K^-$ ; the center-of-mass decay angles are chosen randomly on the unit sphere.

The background particles are generated by the following algorithm:

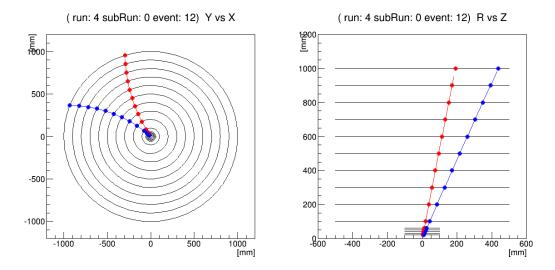
- Background particles are generated in pairs, one  $\pi^+$  and one  $\pi^-$ .
- The number of pairs in each event is a random variate chosen from a Poisson distribution with a mean of 0.75.
- Each of the pions is generated as follows:
  - It is produced at the origin.
  - It has a momentum that is chosen randomly from a distribution that is uniform between 0 and 800 MeV/c.
  - Its direction is chosen randomly on the unit sphere.

The above algorithm generates events with a total charge of zero but there is no concept of momentum or energy balance. About 47% of these events will not have any background tracks.

In the detector simulation step, particles neither scatter nor lose energy when they pass through the detector cylinders; nor do they decay. Therefore, the charged particles follow a perfectly helical trajectory. The simulation follows each charged particle until it either exits the detector or until it completes the outward-going arc of the helix. When the simulated trajectory crosses one of the detector shells, the simulation records the true point of intersection. All intersections are recorded; at this stage in the simulation, there is no notion of inefficiency or resolution. The simulation does not follow the trajectory of the  $\phi$  meson because it was decayed in the generator.

Figure 3.4 shows an event display of a simulated event that has no background tracks. In this event the  $\phi$  meson was travelling close to 90° to the z axis and it decayed nearly symmetrically; both tracks intersect all 15 detector cylinders. The left-hand figure shows an xy view of the event; the solid lines show the trajectory of the kaons, red for  $K^+$  and blue for  $K^-$ ; the solid dots mark the intersections of the trajectories with the detector shells. The right-hand figure shows the same event but in an rz view.

Figure 3.5 shows an event display of another simulated event, one that has four background tracks, all drawn in green. In the xy view it is difficult to see the two  $\pi^-$  tracks, which



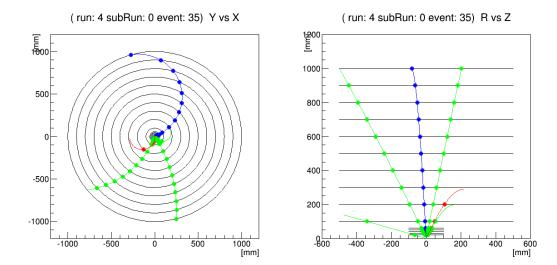
**Figure 3.4:** Event display of a simulated event in the toy detector.

have very low transverse momentum, but they are clear in the rz view. Look at the  $K^+$  track, draw in red; its trajectory just stops in the middle of the detector. Why does this happen? In order to keep the exercises focused on art details, not geometric corner cases, the simulation stops a particle when it completes the outward-going arc of the helix and starts to curl back towards the z axis; it does this even if the particle is still inside the detector.

The third step in the simulation chain (hit-making) is to inspect the intersections produced by the detector simulation and turn them into data-like hits. In this step, a simple model of inefficiency is applied and some intersections will not produce hits. Each hit represents a 2D measurement  $(\varphi, z)$ ; each component is smeared with a gaussian distribution.

The three simulation steps use tools provided by art to record the  $truth\ information(\gamma)$  about each hit. Therefore it is possible to navigate from any hit back to the intersection from which it is derived, and from there back to the particle that made the intersection.

The fourth step is the reconstruction step. The toyExperiment does not yet have properly working reconstruction code; instead it mocks up credible looking results. The output of this code is a data product that represents a fitted helix; it contains the fitted track parameters of the helix, their covariance matrix and collection of smart pointers that point to the

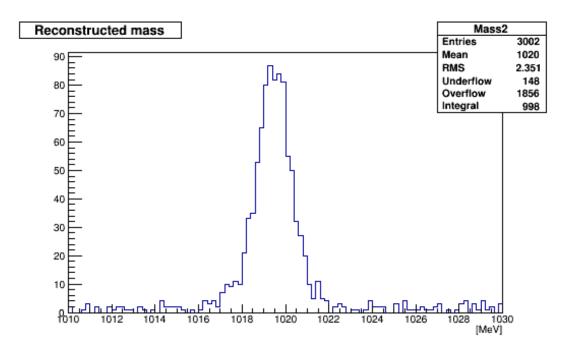


**Figure 3.5:** Event display of another simulated event in the toy detector; a  $K^-$  (blue) is produced with a very shallow trajectory and it does not intersect any detector shells while the  $K^+$  (red) makes five hits in the inner detector and seven in the outer detector

hits that are on the reconstructed track. When we write proper tracking finding and track fitting code for the toyExperiment, the classes that describe the fitted helix will not change. Because the main point of the Workbook exercises is to illustrate the bookkeeping features in *art*, this is good enough for the task at hand. The mocked-up reconstruction code will only create a fitted helix object if the number of hits on a track is greater than some minimum value. Therefore there may be some events in which the output data product is be empty.

The fifth step in the workflow does a simulated analysis using the fitted helices from the reconstruction step. It forms all distinct pairs of tracks and requires that they be oppositely charged. It then computes the invariant mass of the pair, under the assumption that both fitted helices are kaons.\*\* This module is an analyzer module and does not make any output data product. But it does make some histograms, one of which is a histogram of the reconstructed invariant mass of all pairs of oppositely charged tracks; this histogram is shown in Figure 3.6. When you run the Workbook exercises, you will make this plot and can compare it to Figure 3.6. In the figure you can see a clear peak that is created when

<sup>\*\*</sup>The toy experiment does not have any particle identification system so analysis code cannot know if a reconstructed track is a pion or a kaon. A planned enhancement of the toy experiment is to add a particle identification device.



**Figure 3.6:** The invariant mass of all reconstruct pairs of oppositely charged tracks; for this all reconstructed tracks are assumed to be kaons.

the two reconstructed tracks are the two true daughters of the generated  $\varphi$  meson. You can also see an almost flat contribution that occurs when at least one of the reconstructed tracks comes from one of the generated background particles.

## 3.8 Rules, Best Practices, Conventions and Style

In many places, the Workbook will recommend that you write fragments of code in a particular way. The reason for any particular recommendation may be one of the following:

- It is a hard rule enforced by the C++ language or by one of the external products.
- It is a recommended best practice that might not save you time or effort now but will in the long run.
- It is a convention that is widely adopted; C++ is a rich enough language that it will let you do some things in many different ways. Code is much easier to understand

and debug if an experiment chooses to always write code fragments with similar intent using a common set of conventions.

• It is simply a question of style.

It is important to be able to distinguish between rules, best practices, conventions and styles; you must follow the rules; it wise to use best practices and established conventions; but style suggestions are just that, suggestions. This documentation will distinguish among these options when discussing the recommendations that it makes.

If you follow the recommendations for best practices and common conventions, it will be easier to verify that your code is correct and your code will be easier to understand, develop and maintain.

## 4 Unix Prerequisites

#### 4.1 Introduction

You will work through the Workbook exercises on a computer that is running some version of the Unix operating system. This chapter describes where to find information about Unix and gives a list of Unix commands that you should understand before starting the Workbook exercises. This chapter also describes a few ideas that you will need immediately but which are usually not covered in the early chapters of standard Unix references.

If you are already familiar with Unix and the  $bash(\gamma)$  shell, you can safely skip this chapter.

### 4.2 Commands

In the Workbook exercises, most of the commands you will enter at the Unix prompt will be standard Unix commands, but some will be defined by the software tools that are used to support the Workbook. The non-standard commands will be explained as they are encountered. To understand the standard Unix commands, any standard Linux or Unix reference will do. Section 4.10 provides links to Unix references.

Most Unix commands are documented via the *man page* system (short for "manual"). To get help on a particular command, type the following at the command prompt, replacing <command-name> with the actual name of the command:

#### man <command-name>

In Unix, everything is case sensitive; so the command man must be typed in lower case. You can also try the following; it works on some commands and not others:

```
<command-name> -help
```

or

#### <command-name> -?

Before starting the Workbook, make sure that you understand the basic usage of the following Unix commands:

```
cat, cd, cp, echo, export, gzip, head, less, ln -s, ls, mkdir, more, mv, printenv, pwd, rm, rmdir, tail, tar
```

You also need to be familiar with the following Unix concepts:

- o filename vs pathname
- o absolute path vs relative path
- o directories and subdirectories (equivalent to folders in the Windows and Mac worlds)
- o current working directory
- home directory (aka login directory)
- . . / notation for viewing the directory above your current working directory
- environment variables (discussed briefly in Section 4.5)
- $\circ$  paths( $\gamma$ ) (in multiple senses; see Section 4.6)
- file protections (read-write-execute, owner-group-other)
- o symbolic links
- o stdin, stdout and stderr
- redirecting stdin, stdout and stderr
- o putting a command in the background via the & character
- o pipes

#### 4.3 Shells

When you type a command at the prompt, a Unix agent called a *Unix shell*, or simply a *shell*, reads your command and figures out what to do. Some commands are executed internally by the shell but other commands are dispatched to an appropriate program or script. A shell lives between you and the underlying operating system; most versions of Unix support several shells. The *art* Workbook code expects to be run in the *bash shell*. You can see which shell you're running by entering:

#### echo \$SHELL



For those of you with accounts on a Fermilab machine, your login shell was initially set to the *bash* shell\*.



If you are working on a non-Fermilab machine and *bash* is not your default shell, consult a local expert to learn how to change your login shell to bash.

## 4.4 Scripts: Part 1

In order to automate repeated operations, you may write multiple Unix commands into a file and tell *bash* to run all of the commands in the file as if you had typed them sequentially. Such a file is an example of a *shell script* or a *bash script*. The *bash* scripting language is a powerful language that supports looping, conditional execution, tests to learn about properties of files and many other features.

Throughout the Workbook exercises you will run many scripts. You should understand the big picture of what they do, but you don't need to understand the details of how they work.



If you would like to learn more about bash, some references are listed in Section 4.10.

<sup>\*</sup> If you have had a Fermilab account for many years, your default shell might be something else. If your default shell is not bash, open a Service Desk ticket to request that your default shell be changed to bash.

#### 4.5 Unix Environments

#### 4.5.1 Building up the Environment

Very generally, a Unix *environment* is a set of information that is made available to programs so that they can find everything they need in order to run properly. The Unix operating system itself defines a generic environment, but often this is insufficient for everyday use. However, an environment sufficient to run a particular set of applications doesn't just pop out of the ether, it must be *established* or *set up*, either manually or via a script. Typically, on institutional machines at least, system administrators provide a set of login scripts that run automatically and enhance the generic Unix environment. This gives users access to a variety of system resources, including, for example:

- disk space to which you have read access
- o disk space to which you have write access
- o commands, scripts and programs that you are authorized to run
- o proxies and tickets that authorize you to use resources available over the network
- the actual network resources that you are authorized to use, e.g., tape drives and DVD drives

This constitutes a basic *working environment* or *computing environment*. Environment information is largely conveyed by means of *environment variables* that point to various program executable locations, data files, and so on. A simple example of an environment variable is HOME, the variable whose value is the absolute path to your home directory.

Particular programs (e.g., *art*) usually require extra information, e.g., paths to the program's executable(s) and to its dependent programs, paths indicating where it can find input files and where to direct its output, and so on. In addition to environment variables, the *art*-enabled computing environment includes some aliases and *bash* functions that have been defined; these are discussed in Section 4.8.

In turn, the Workbook code, which must work for all experiments and at Fermilab as well as at collaborating institutions, requires yet more environment configuration – a *site-specific* configuration.

Given the different experiments using art and the variety of laboratories and universities at which the users work, a  $site(\gamma)$  in art is a unique combination of experiment and institution. It is used to refer to a set of computing resources configured for use by a particular experiment at a particular institution. Setting up your site-specific environment will be discussed in Section 4.7.



When you finish the Workbook and start to run real code, you will set up your experiment-specific environment on top of the more generic *art*-enabled environment, in place of the Workbook's. To switch between these two environments, you will log out and log back in, then run the script appropriate for the environment you want. Because of potential naming "collisions," it is not guaranteed that these two environments can be overlain and always work properly.

This concept of the environment hierarchy is illustrated in Figure 4.1.

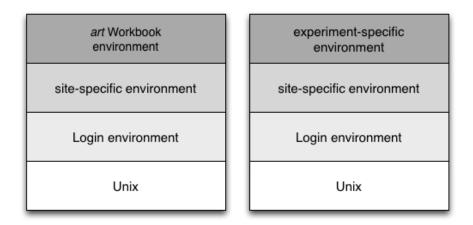


Figure 4.1: Hierarchies of the *art* Workbook (left) and experiment-specific (right) computing environments

### 4.5.2 Examining and Using Environment Variables

One way to see the value of an environment variable is to use the printenv command:

### printenv HOME

At any point in an interactive command or in a shell script, you can tell the shell that

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you want the value of the environment variable by prefixing its name with the \$ character:

#### echo \$HOME

Here, echo is a standard Unix command that copies its arguments to its output, in this case the screen.

By convention, environment variables are virtually always written in all capital letters<sup>†</sup>.

There may be times when the Workbook instructions tell you to set an environment variable to some value. To do so, type the following at the command prompt:

```
export <ENVNAME>=<value>
```

If you read *bash* scripts written by others, you may see the following variant, which accomplishes the same thing:

```
<ENVNAME>=<value>export <ENVNAME>
```

## **4.6 Paths and \$PATH**

*Path* (or *PATH*) is an overloaded word in computing. Here are the ways in which it is used:

**Lowercase** path can refer to the location of a file or a directory; a path may be absolute or relative, e.g.

```
/absolute/path/to/mydir/myfile or relative/path/on/same/branch/to/mydir/myfile or ../relative/path/on/different/branch/to/herdir/herfile
```

PATH refers to the standard Unix environment variable set by your login scripts and updated by other scripts that extend your environment; it is a colon-separated list of directory names, e.g.,

```
/usr/bin:/usr/sbin:/usr/local/bin.
```

<sup>&</sup>lt;sup>†</sup>Another type of variable, *shell variables*, are local to the currently-invoked shell and go away when the shell exits. By convention, these are written in lower or mixed case. These conventions provide a clue to the programmer as to whether changing a variable's value might have consequences outside the current shell.

It contains the list of directories that the shell searches to find programs/files required by Unix shell commands (i.e., PATH is used by the shell to "resolve" commands).

"path" generically, refers to any environment variable whose value is a colon-separated list of directory names e.g.,

```
/abs/path/a:/abs/path/b:rel/path/c
```

In addition, *art* defines a fourth idea, also called a path, that is unrelated to any of the above; it will be described as you encounter it in the Workbook, e.g., Section 9.8.8.

All of these path concepts are important to users of *art*. In addition to PATH itself, there are three PATH-like environment variables (colon-separated list of directory names) that are particularly important:

```
LD_LIBRARY_PATH used by art to resolve shareable libraries
PRODUCTS used by UPS to resolve external products
```

FHICL\_FILE\_PATH use by FHiCL to resolve #include directives.

When you source the scripts that setup your environment for *art*, these will be defined and additional colon-separated elements will be added to your PATH. To look at the value of PATH (or the others), enter:

```
printenv PATH
```

To make the output easier to read by replacing all of the colons with newline characters, enter:

```
printenv PATH | tr : \\n
```

In the above line, the vertical bar is referred to as a *pipe* and tr is a standard Unix command. A pipe takes the output of the command to its left and makes that the input of the command to its right. The tr command replaces patterns of characters with other patterns of characters; in this case it replaces every occurrence of the colon character with the new-line character. To learn why a double back slash is needed, read bash documentation to learn about escaping special characters.

## 4.7 Scripts: Part 2

There are two ways to run a bash script (actually three, but two of them are the same). Suppose that you are given a bash script named file.sh. You can run any of these commands:

file.sh

source file.sh

. file.sh

The first version, file.sh, starts a new bash shell, called a subshell, and it executes the commands from file.sh in that subshell; upon completion of the script, control returns to the parent shell. At the startup of a subshell, the environment of that subshell is initialized to be a copy of the environment of its parent shell. If file.sh modifies its environment, then it will modify only the environment of the subshell, leaving the environment of the parent shell unchanged. This version is called *executing* the script.

The second and third versions are equivalent. They do not start a subshell; they execute the commands from file.sh in your current shell. If file.sh modifies any environment variables, then those modifications remain in effect when the script completes and control returns to the command prompt. This is called *sourcing* the script.

Some shell scripts are designed so that they must be sourced and others are designed so that they must be executed. Many shell scripts will work either way.



If the purpose of a shell script is to modify your working environment then it must be sourced, not executed. As you work through the Workbook exercises, pay careful attention to which scripts it tells you to source and which to execute. In particular, the scripts to setup your environment (the first scripts you will run) are bash scripts that must be sourced because their purpose is to configure your environment so that it is ready to run the Workbook exercises.

Some people adopt the convention that all bash scripts end in .sh; others adopt the convention that only scripts designed to be sourced end in .sh while scripts that must be executed have no file-type ending (no ".something" at the end). Neither convention is uniformly applied either in the Workbook or in HEP in general.

If you would like to learn more about bash, some references are listed in Section 4.10.

## 4.8 bash Functions and Aliases

The bash shell also has the notion of a *bash function*. Typically bash functions are defined by sourcing a bash script; once defined, they become part of your environment and they can be invoked as if they were regular commands. The setup product "command" that you will sometimes need to issue, described in Chapter 7, is an example. A bash function is similar to a bash script in that it is just a collection of bash commands that are accessible via a name; the difference is that bash holds the definition of a function as part of the environment while it must open a file every time that a bash script is invoked.

You can see the names of all defined bash functions using:

#### declare -F

The bash shell also supports the idea of *aliases*; this allows you to define a new command in terms of other commands. You can see the definition of all aliases using:

#### alias

You can read more about bash shell functions and aliases in any standard bash reference.

When you type a command at the command prompt, bash will resolve the command using the following order:

- 1. Is the command a known alias?
- 2. Is the command a bash keyword, such as if or declare?
- 3. Is the command a shell function?
- 4. Is the command a shell built-in command?
- 5. Is the command found in \$PATH?

To learn how bash will resolve a particular command, enter:

type <command-name>

## 4.9 Login Scripts

When you first login to a computer running the Unix operating system, the system will look for specially named files in your home directory that are scripts to set up your working environment; if it finds these files it will source them before you first get a shell prompt. As mentioned in Section 4.5, these scripts modify your PATH and define bash functions, aliases and environment variables. All of these become part of your environment.



When your account on a Fermilab computer was first created, you were given standard versions of the files .profile and .bashrc; these files are used by bash<sup>‡</sup>. You can read about login scripts in any standard bash reference. You may add to these files but you should not remove anything that is present.



If you are working on a non-Fermilab computer, inspect the login scripts to understand what they do.

It can be useful to inspect the login scripts of your colleagues to find useful customizations.

If you read generic Unix documentation, you will see that there are other login scripts with names like, .login, .cshrc and .tcshrc. These are used by the csh family of shells and are not relevant for the Workbook exercises, which require the bash shell.

## 4.10 Suggested Unix and bash References

The following cheat sheet provides some of the basics:

http://mu2e.fnal.gov/atwork/computing/UnixHints.shtml

A more comprehensive summary is available from:

o http://www.tldp.org/LDP/.../GNU-Linux-Tools-Summary.html

Information about writing bash scripts and using bash interactive features can be found in:

 BASH Programming - Introduction HOW-TO http://tldp.org/HOWTO/Bash-Prog-Intro-HOWTO.html

<sup>&</sup>lt;sup>‡</sup>These files are used by the sh family of shells, which includes bash.

- Bash Guide for Beginners
   http://www.tldp.org/LDP/Bash-Beginners-Guide/html/Bash-Beginners-Guide.html
- Advanced Bash Scripting Guide http://www.tldp.org/LDP/abs/html/abs-guide.html

The first of these is a compact introduction and the second is more comprehensive.

The above guides were all found at the Linux Documentation Project: Workbook:

http://www.tldp.org/guides.html

# 5 Site-Specific Setup Procedure

Section 4.5 discussed the notion of a working environment on a computer. This chapter answers the question: How do I make sure that my environment variables are set correctly to run the Workbook exercises or my experiment's code using *art*?

Very simply, on every computer that hosts the Workbook, a procedure must be established that every user is expected to follow once per login session. In most cases (NO $\nu$ A being a notable exception), the procedure involves only sourcing a shell script (recall the discussion in Section 4.7). In this documentation, we refer to this procedure as the "site-specific setup procedure." It is the responsibility of the people who maintain the Workbook software for each  $site(\gamma)$  to ensure that this procedure does the right thing on all the site's machines.



As a user of the Workbook, you will need to know what the procedure is and you must remember to follow it each time that you log in.



For all of the Intensity Frontier experiments at Fermilab, the site-specific setup procedure defines all of the environment variables that are necessary to create the working environment for either the Workbook exercises or for the experiment's own code.

Table 5.1 lists the site-specific setup procedure for each experiment. You will follow the procedure when you get to Section 9.6.

**Table 5.1:** Site-specific setup procedures for  $IF(\gamma)$  Experiments at Fermilab; for the equivalent procedure at a non-Fermi site, consult an expert from that site. For the NO $\nu$ A experiment, the procedure is a sequence of three commands. For all others it is a single command, but in a few cases the command is so long that the table shows it split over two lines: you must type it as a single command on one line; the experiments whose command is so split are ArgoNeut, LBNE and MicroBoone.

Experiment	Site-Specific Setup Command
ArgoNeut	source /grid/fermiapp/argoneut/code/t962soft/setup/
	setup_t962_fnal.sh -r art_workbook -b prof
Darkside	source /ds50/app/ds50/ds50.sh
LBNE	<pre>source /grid/fermiapp/lbne/lar/code/larsoft/setup/</pre>
	setup_larsoft_fnal.sh -r art_workbook -b prof
MicroBoone	<pre>source /grid/fermiapp/lbne/lar/code/larsoft/setup/</pre>
	setup_larsoft_fnal.sh -r art_workbook -b prof
Muon g-2	source /gm2/app/software/prod/g-2/setup
Mu2e	setup mu2e
ΝΟνΑ	<pre>source /grid/fermiapp/nova/novaart/novasvn/srt/srt.sh</pre>
	export EXTERNALS=/nusoft/app/externals
	<pre>source \$SRT_DIST/setup/setup_novasoft.sh -b maxopt</pre>

# 6 Get your C++ up to Speed

## 6.1 Introduction

,

There are two goals for this chapter. The first is to illustrate the features of C++ that will be important for users of the Workbook, especially those features that will be used in the first few Workbook exercises. It does not attempt to cover C++ comprehensively and it delegates as much as possible to the standard documentation.

The second goal is to explain the process of turning source code files into an *executable program*. The two steps in this process are *compiling* and *linking*. In informal writing, the word *build* is sometimes used to mean just compiling or just linking, but usually it refers to the two together.

A typical program consists of many source code files, each of which contains a human-readable description of one component of the program. In the Workbook, you will see source code files written in the C++ computer language; these files have names that end in .cc. In C++, there is a second sort of source code file, called a *header file* that ends in .h; in most, but not all, cases for every file ending in .cc there is another file with the same name but ending in .h. Header files can be thought of as the "parts list" for the corresponding .cc file; you will see how these are used in Section 6.4.

In the compilation step each .cc file is translated into *machine code*, also called *binary code* or *object code*, which is a set of instructions, in the computer's native language, to do the tasks described by the source code. The output of the compilation step is called an *object file*; in the examples you will see in the Workbook, object files always end in .o. But an object file, by itself, is not an *executable program*. It is not executable because each

. o file was created in isolation and does not know about the other . o files.

It is often convenient to collect related groups of .o files and put them into *libraries*. There are two kinds of library files, static libraries, whose names end in .a and shared libraries whose names end in .so. Putting many .o files into a single library allows you to use them as a single coherent entity. We will defer further discussion of libraries until more background information has been provided.

The job of the *linking* step is to read the information found in the various libraries and .o files and form them into an *executable program*. When you run the linker, you tell it the name of the file into which it will write the executable program. It is a common, but not universal, practice that the filename of an executable program has no extension (i.e., no .something at the end).

After the linker has finished, you can run your executable program typing the filename of the program at the bash command prompt.

A typical program links both to libraries that were built from the program's source code and to libraries from other sources. Some of these other libraries might have been developed by the same programmer as general purpose tools to be used by his or her future programs; other libraries are provided by third parties, such as *art* or your experiment. Many C++ language features are made available to your program by telling the linker to use libraries provided by the C++ compiler vendor. Other libraries are provided by the operating system.

Now that you know about libraries,, we can give a second reason why an object file, by itself, is not an executable program: until it is linked, it does not have access to the functions provided by any of the external libraries. Even the simplest program will need to be linked against some of the libraries supplied by the compiler vendor and by the operating system.

The names of all of the libraries and object files that you give to the linker is called the *link list*.

This chapter is designed around a handful of exercises, each of which you will first build and run, then "pick apart" to understand how the results were obtained.

## 6.2 Establishing the Environment

### 6.2.1 Initial Setup

To start these exercises for the first time, do the following:

- 1. Log into the node that you will use for Workbook exercises.
- 2. Follow the site-specific setup procedure from Table 5.1.
- 3. Create an empty working directory and cd to it.
- 4. Run these commands to copy a gzipped tar file from the web, unpack it, and get a directory listing:

```
wget https://sharepoint.fnal.gov/project/ \ArtDoc-Pub/Shared%20Documents/C++Up
tar xzf C++UpToSpeed.tar.gz
rm C++UpToSpeed.tar.gz
ls
BasicSyntax Build Classes Libraries setup.sh
```

5. To select the correct compiler version and define a few environment variables that will be used later in these exercises, run:

```
source setup.sh
```

After these steps, you are ready to begin the exercise in Section 6.3.

## 6.2.2 Subsequent Logins

If you log out and log back in again, reestablish your environment by following these steps:

- 1. Log into the node that you will normally use.
- 2. Follow the site-specific setup procedure.
- 3. cd to the working directory you created in Section 6.2.1.

- 4. source setup.sh
- 5. cd to the directory that contains the exercise you want to work on.

## 6.3 C++ Exercise 1: The Basics

## 6.3.1 Concepts to Understand

This section provides a program that illustrates the parts of C++ that are assumed knowledge for the Workbook material. If you do not understand some of the code in this example program, consult any standard C++ reference; several are listed in Section 6.7.

Once you have understood this example program, you should understand the following C++ concepts:

- o how comments are indicated
- o what is a main program
- o how to write a main program
- o how to compile, link and run the main program
- o how to distinguish between source, object and executable files
- o how to print to standard output, std::cout
- o how to declare and define  $variables(\gamma)$  of the some of the frequently used built-in types: int, float, double, bool
- the {} initializer syntax
- o assignment to variables
- o arrays
- o several different forms of looping
- o comparisons: ==, !=, <, >, >=, <=
- o if-then-else, if-then-else if-else

- o pointers
- o references
- o std::string (a type from the C++ Standard Library ( $std(\gamma)$ )
- the class template from the standard library, std::vector<T>\*

The above list explicitly does not include classes, objects and inheritance, which will be discussed in Sections 6.6 and 32.9.

## 6.3.2 How to Compile, Link and Run

In this section you will learn how to compile, link and run the small C++ program that illustrates the features of C++ that are considered prerequisites. The main discussion of the details of compiling and linking will be deferred until Section 6.4.

We don't offer a lot of details up front; more will follow in Sections 6.3.5 and 6.3.4. The idea here is to get used to the steps and see what results you get.

To compile, link and run the sample C++ program, called t1:

- 1. If not yet done, log in and establish the working environment (Section 6.2).
- 2. List the starting set of files:

```
cd BasicSyntax/v1/
Is
build t1.cc t1_example.log
```

The file t1.cc contains the source code of the *main pro*gram, which is a function called main() { ...}. The file build is a script that will compile and link the code. The file  $t1\_example.log$  is an example of the output expected when

<sup>\*</sup>You need to know how to use std::vector<T> but you do not need to understand how it works or how to write your own templates.

you run t1.

3. Compile and link the code (one step); then look at a directory listing:

#### build

```
t1.cc: In function 'int main()':
t1.cc:43:26: warning: 'k' may be
used uninitialized in this function
[-Wuninitialized]
```

#### ls

```
build t1 t1.cc t1_example.log
```

The script named build compiles and links the code, and produces the executable file t1. The warning message, issued by the compiler, also comes during this step.

4. Run the executable file sending output to a log file:

```
./t1 > t1.log
```

### 6.3.3 Suggested Homework

1. Compare your output with the standard example:

```
diff t1.log t1_example.log
```

There will almost certainly be a handful of differences.

2. Look at the file t1.cc and understand what it does, in particular the relationship between the lines in the program and the lines in the output.

If you don't understand something, consult a standard C++ reference; see Section 6.7. A few of your questions might also be answered in Section 6.3.4.

#### 6.3.4 Discussion

Why do we expect several of the lines of the output to be different from those in t1\_example.log? There are two classes of answers: (1) an uninitialized variable and (2) variation in variable addresses from run to run.

In t1.cc, the line int k; declares that k is a variable whose type is int but it does not initialize the variable. Therefore the value of the variable k is whatever value happened to be sitting in the memory location that the program assigned to k. Each time that the program runs, the operating system will put the program into whatever region of memory makes sense to the operating system; therefore the address of any variable, and thus the value returned, may change unpredictably from run to run.

This line is also the source of the warning message produced by the build script. This line was included to make it clear what we mean by *initialized* variables and *uninitialized* variables. Uninitialized variables are frequent sources of errors in code and therefore you should *always* initialize your variables. In order to help you establish this good coding habit, the remaining exercises in this series and in the Workbook include the compiler option <code>-Werror</code>. This tells the compiler to promote warning messages to error level and to stop compilation without producing an output file.

The second line that may differ from one run to the next is:

```
float *pa=&a;
```

This line declares a variable pa, which is of type  $pointer(\gamma)$  to float, and it initializes it to be the memory address of the variable a (a must be of type float). Since the address may change from run to run, so may the printout that starts pa =.

For similar reasons, the lines in the printout that start &a = and &ra = may also change from run to run.

## 6.3.5 How was this Exercise Built?

Just to see how the exercise was built, look at the script BasicSyntax/v1/build that you ran to compile and link t1.cc; the following command was issued:

```
c++ -Wall -Wextra -pedantic -std=c++11 -o t1 t1.cc
```

Part

This turned the source file t1.cc into an executable program, named t1 (the argument to the -o (for "output") option). We will discuss compiling and linking in Section 6.4.

## 6.4 C++ Exercise 2: About Compiling and Linking

#### 6.4.1 What You Will Learn

In the previous exercise, the entire program was found in a single file and the build script performed compiling and linking in a single step. For all but the smallest programs, this is not practical. It would mean, for example, that you would need to recompile and relink everything when you made even the smallest change anywhere in the code; generally this would take much too long. To address this, some computer languages, including C++, allow you to break up a large program into many smaller files and rebuild only a small subset of files when you make changes in one.

There are two exercises in this section. In the first one the source code consists of three files. This example has enough richness to discuss the details of what happens during compiling and linking, without being overwhelming. The second exercise introduces the ideas of libraries and external packages.

#### 6.4.2 The Source Code for this Exercise

The source code for this exercise is found in Build/v1, relative to your working directory. The relevant files are function.cc, function.h and t1.cc.

The file t1.cc contains the source code for the function main() { ...} for this exercise. Every C++ program must have one and only one function named main, which is where the program actually starts execution. Note that the term *main program* sometimes refers to this function, but other times refers to the .cc file that contains it. In either case, *main program* refers to this function, either directly or indirectly. For more information, consult any standard C++ reference. The file function.h is a header file that declares a function named function. The file function.cc is another source code file; it provides the definition of that function.



Look at t1.cc: it both declares and defines the program's function main() { ...} that takes no arguments. A function with this  $signature(\gamma)$  has special meaning to the com-

plier and the linker: they recognize it as a C++ main program. There are other signatures that the compiler and linker will recognize as a C++ main program; consult the standard C++ documentation.



To be recognized as a main program, there is one more requirement: main() { ...} must be declared in the global namespace.

The body of the main program (between the braces), declares and defines a variable a and initializes it to the value of 3; it prints out the value of a. Then it calls a function that takes a as an argument and prints out the value returned by that function.

You, as the programmer using that function, need to know what the function does but the C++ compiler doesn't. It only needs to know the name, argument list and return type of the function — information that is provided in the header file, function.h. This file contains the line

```
float function( float );
```

This line is called the  $declaration(\gamma)$  of the function. It says (1) that the identifier function is the name of a function that (2) takes an argument of type float (the "float" inside the parentheses) and (3) returns a value of type float (the "float" at the start of the line). The file t1.cc includes this header file, thereby giving the compiler these three pieces of information it needs to know about function.

The other three lines in function.h are *code guards*, described in Section 32.8. In brief, they deal with the following scenario: suppose that we have two header files, A.h and B.h, and that A.h includes B.h; there are many scenarios in which it makes good sense for a third file, either .h or .cc, to include both A.h and B.h. The code guards ensure that, when all of the includes have been expanded, the compiler sees exactly one copy of B.h.

Finally, the file function.cc contains the source code for the function named function:

```
float function ( float i ) {
  return 2.*i;
}
```

It names its argument i, multiplies this argument by two and returns that value. This code fragment is called the *definition* of the function or the *implementation*( $\gamma$ ) of the function.

(The C++ standard uses the word *definition* but *implementation* is in common use.)

We now have a rich enough example to discuss another case in which the same word is frequently used to mean two different things. Sometimes people use the phrase "the source code of the function named function" to refer collectively to both function.h and function.cc; sometimes they use it to refer exclusively to function.cc. Unfortunately the only way to distinguish the two uses is from context.

The word *header file* always refers unambiguously to the .h file. The term *implementation file* is used to refer unambiguously to the .cc file. This name follows from the its contents: it describes how to implement the items declared in the header file.

Based on the above description, when this exercise is run, we expect it to print out:

```
a = 3 function(a) 6
```

## 6.4.3 Compile, Link and Run the Exercise

To perform this exercise, first log in and cd to your working directory if you haven't already, then

1. cd to the directory for this exercise and get a directory listing:

```
cd Build/v1
ls
build build2 function.cc function.h t1.cc
```

The two files, build and build2 are scripts that show two different ways to build the code.

2. Compile and link this exercise, then get an updated directory listing:

```
build
ls
build build2 function.cc function.h
```

```
function.o t1 t1.cc t1.o

Notice the new files function.o, t1 and t1.o.

3. Run the exercise:

./t1
a = 3
function(a) 6
```

This matches the expected printout.

Look at the file build that you just ran. It has three steps; the first two commands have the -c command line option while the last one does not:

- 1. It compiles the main program, t1.cc, into the object file (with the default name) t1.o (which will now be the thing that the term *main program* refers to): c++ -Wall -Wextra -pedantic -Werror -std=c++11 -c t1.cc
- 2. It (separately) compiles function.cc into the object file function.o (shown in two lines):

```
c++ -Wall -Wextra -pedantic -Werror -std=c++11 -c function.cc
```

3. It links t1.0 and function.0 to form the executable program t1 (the name of the main program is the argument of the -o option):

```
c++ -std=c++11 -o t1 t1.o function.o
```

You should have noticed that the same command, c++, is used both for compiling and linking. The full story is that when you run the command c++, you are actually running a program that parses its command line to determine which, if any, files need to be compiled and which, if any, files need to be linked. It also determines which of its command line arguments should be forwarded to the compiler and which to the linker. It then runs the compiler and linker as many times as required.

If the -c option is present, it tells c++ to compile only, and not to link. If -c is specified, the .cc file(s) to compile must also be specified. Each of the files will be compiled to

create its corresponding object file and then processing stops. In our example, the first two commands each compile a single source file. Note that if any .o files are given on the command line, c++ will issue a warning and ignore them.

The third command (with no -c option) is the linking step. Even if the -c option is missing, c++ will first look for source files on the command line; if it finds any, it will compile them and put the output into temporary object files. In our example, there are none, so it goes straight to linking. The two just-created object files are specified (at the end, here, but the order is not important); the -o t1 portion of the command tells the linker to write its output (the executable) to the file t1.

As it is compiling the main program, t1.cc, the compiler recognizes every function that is defined within the file and every function that is called by the code in the file. It recognizes that t1.cc defines a function main() and that main() calls a function named function, whose definition is not found inside t1.cc. At the point that t1.cc calls function, the compiler will write to function all of the machine code needed to prepare for the call; it will also write all of the machine code needed to use the result of the function. In between these two pieces, the compiler will write machine code that says "call the function whose memory address is" but it must leave an empty placeholder for the address. The placeholder is empty because the compiler does not know the memory address of that function.

The compiler also makes a table that lists all functions defined by the file and all functions that are called by code within the file. The name of each entry in the table is called a *linker symbol* and the table is called a *symbol table*. When the compiler was compiling t1.cc and it found the definition of the main program, it created a linker symbol for the main program and added a notation to say the this file contains the definition of that symbol. When the compiler was compiling t1.cc and it encountered the call to function, it created a linker symbol for this function; it marked this symbol as an *undefined reference* (because it could not find the definition of function within t1.cc). The symbol table also lists all of the places in the machine code of t1.o that are placeholders that must be updated once the memory address of function is known. In this example there is only one such place.

When the compiler writes an object file, it writes out both the compiled code and the table of linker symbols.

In t1.cc, the compiled code for the line that begins std::cout will do its work by calling a few functions that are found in the compiler-supplied libraries. The linker symbols for these functions will also be listed as undefined references in the symbol table of t1.o; the symbol table also lists the places within the machine code of t1.o that need to be updated once the addresses of these symbols are known.

The symbol table in the file function o is simple; it says that this file defines a function named function that takes a single argument of type float and that returns a float.

The job of the linker (also invoked by the command c++) is to play match-maker. First it inspects the symbol tables inside all of the object files listed on the command line and looks for a linker symbol that defines the location of the main program. If it cannot find one, or if it finds more than one, it will issue an error message and stop. In this example

- 1. The linker will find the definition of a main program in t1.0.
- 2. It will start to build the executable (output) file by copying the machine code from t1.0 to the output file.
- 3. Then it will try to resolve the unresolved references listed in the symbol table of t1.0; it does this by looking at the symbol tables of the other object files on the command line. It also knows to look at the symbol tables from a standard set of compiler-supplied and system-supplied libraries.
- 4. It will discover that function o resolves one of the external references from t1.0. So it will copy the machine code from function o to the executable file.
- 5. It will discover that the other unresolved references in t1.0 are found in the compiler-supplied libraries and will copy code from these libraries into the executable.
- 6. Once all of the machine code has been copied into the executable, the compiler knows the memory address of every function. The compiler can then go into the machine code, find all of the placeholders and update them with the correct memory addresses.

Sometimes resolving one unresolved reference will generate new ones. The linker iterates until (a) all references are resolved and no new unresolved references appear (success)

or (b) the same unresolved references continue to appear (error). In the former case, the linker writes the output to the file specified by the -o option; if no -o option is specified the linker will write its output to a file named a .out. In the latter case, the linker issues an error message and does not write the output file.

After the link completes, the files t1.0 and function.0 are no longer needed because everything that was useful from them was copied into the executable t1. You may delete the .0 files, and the executable will still run.

## **6.4.4** Alternate Script build2

The script build2 shows an equivalent way of building t1 that is commonly used for small programs; it does it all on one line. To exercise this script:

- 1. Stay in the same directory as before, Build/v1.
- 2. Clean up from the previous build and look at the directory contents:

```
rm function.o t1 t1.o
ls
build build2 function.cc function.h t1.cc
```

3. Run the build2 script, and again look at directory contents:

```
build2
ls
build build2 function.cc function.h t1 t1.cc
```

Note that t1 was created but there are no .0 files.

4. Execute the program that you just built:

```
\frac{./t1}{a} = 3
```

```
function(a) 6
```

Look at the script build2; it contains only one command:

```
c++ -Wall -Wextra -pedantic -Werror -std=c++11 -o t1 t1.cc function.cc
```

This script automatically does the same operations as build but it knows that the .o files are temporaries. Perhaps the command c++ kept the contents of the two .o files in memory and never actually wrote them out as disk files. Or, perhaps, the command c++ did explcitly create disk files and deleted them when it was finished. In either case you don't see them when you use build2.

## 6.4.5 Suggested Homework

It takes a bit of experience to decipher the error messages issued by a C++ compiler. The three exercises in this section are intended to introduce you to them so that you (a) get used to looking at them and (b) understand these particular errors if/when you encounter them later.

Each of the following three exercises is independent of the others. Therefore, when you finish with each exercise, you will need to undo the changes you made in the source file(s) before beginning the next exercise.

- 1. In Build/v1/t1.cc, comment out the include directive for function.h; rebuild and observe the error message.
- 2. In Build/v1/function.cc, change the return type to double; rebuild and observe the error message.
- 3. In Build/v1/t1.cc, change float a=3. to double a=3.; rebuild and run. This will work without error and will produce the same output as before.

The first homework exercise will issue the diagnostic:

```
t1.cc: In function 'int main()':
t1.cc:10:44: error: 'function' was not declared in this scope
```

When you see a message like this one, you can guess that either you have not included a required header file or you have misspelled the name of the function.

The second homework exercise will issue the diagnostic (second and last lines split into two here):

```
function.cc: In function 'double function(float)':
function.cc:3:27: error: new declaration
'double function(float)'
In file included from function.cc:1:0:
function.h:4:7: error: ambiguates old declaration
'float function(float)'
```

This error message says that the compiler has found two functions that have the same signature but different return types. The compiler does not know which of the two functions you want it to use.

The bottom line here is that you must ensure that the definition of a function is consistent with its declaration; and you must ensure that the use of a function is consistent with its declaration.

The third homework exercise illustrates the C++ idea of *automatic type conversion*; in this case the compiler will make a temporary variable of type float and set its value to that of a:

```
float tmp = a;
```

The compiler will then use this temporary variable as the argument of the function. Consult the standard C++ documentation to understand when automatic type conversions may occur; see Section 6.7.

## 6.5 C++ Exercise 3: Libraries

Multiple compiled object code files can be grouped into a single file known as a *library*, obviating the need to specify each and every object file when linking; you can reference the libraries instead. This simplifies the multiple use and sharing of software components.

Two Linux C/C++ library types can be created:

o static libraries of object code (filenames for which end in .a) that are linked with, and become part of, the application (*art* does not use static libraries)

dynamically linked, shared object libraries (filenames end in .so); multiple art jobs running simultaneously can dynamically load and unload the same copy of a library of this kind instead of making an exclusive copy of it; this substantially reduces the amount of memory needed by each job.

#### 6.5.1 What You Will Learn

In this section you will repeat the example of Section 6.4 with a variation. You will create an object library, insert function. o into that library and use that library in the link step. This pattern generalizes easily to the case that you will encounter in your experiment software, where object libraries will typically contain many object files.

## 6.5.2 Building and Running the Exercise

To perform this exercise, do the following:

- 1. Log in and establish your working environment (Section 6.2).
- 2. cd to your working directory.
- 3. cd to the directory for this exercise and get a directory listing:

```
cd Libraries/v1
```

ls

build build2 build3 function.cc function.h
t1.cc

The three files, function.cc, function.h and t1.cc are identical to those from the previous exercise. The three files, build, build2 and build3 are scripts that show three different ways to build the main program in this exercise.

4. Compile and link this exercise using build, then compare the

### directory listing to that taken pre-build:

```
build
ls
build build3 function.h libpackage1.a t1.cc
build2 function.cc function.o t1 t1.o
```

5. Execute the main program:

```
./t1
a = 3
function(a) 6
```

This matches the expected printout. Now let's look at the script build. It has four parts which do the following things:

- 1. Compiles function.cc; the same as the previous exercise: c++ -Wall -Wextra -pedantic -Werror -std=c++11 -c function.cc
- 2. Creates the library named libpackage1.a and adds function.o to it: ar rc libpackage1.a function.o

  Note that the name of the library must come before the name of the object file.
- 3. Compiles t1.cc; the same as the previous exercise: c++ -Wall -Wextra -pedantic -Werror -std=c++11 -c t1.cc
- 4. Links the main program against libpackage1.a and the system libraries: c++ -o t1 t1.o libpackage1.a

The two new features are in step 2, which creates the object library, and step 4, in which function o is replaced in the link list with libpackagel.a. If you have many of files to add to the library, you may add them one at a time by repeating step 2 or you may add them all in one command. When you do the latter you may name each object file separately or may use a wildcard:

```
ar rc libpackage1.a *.o
```

In libpackage1.a the string package1 has no special meaning; it was an arbitrary name chosen for this exercise. Actually it was chosen in anticipation of a future exercise that is not yet written up.



The other parts of the name, the prefix lib and the suffix .a, are part of a long-standing Unix convention and some Unix tools presume that object libraries are named following this convention. You should always follow this convention. The use of this convention is illustrated by the scripts build2 and build3.

To perform the exercise using build2, stay in the same directory and cleanup then rebuild as follows:

1. Remove files built by build1:

```
rm function.o t1.o libpackage1.a t1
```

2. Build the code with build2 and look at the directory contents:

```
build2
```

ls

build build3 function.h libpackage1.a t1.cc build2 function.cc function.o t1 t1.o

3. Run ./t1 as before.

The only difference between build and build2 is the link line. The version from build is:

```
c++ -o t1 t1.o libpackage1.a
```

while that from build2 is:

```
c++ -o t1 t1.o -L. -lpackage1
```

In the script build, the path to the library, relative or absolute, is written explicitly on the command line. In the script build2, two new elements are introduced. The command line may contain any number of -L options; the argument of each option is the name of a directory. The ensemble of all of the -L options forms a search path to look for named libraries; the path is searched in the order in which the -L options appear on the line. The names of libraries are specified with the -l options (this is a lower case letter L, not the numeral one); if a -l option has an argument of XXX (or package1), then the linker with search the path defined by the -L options for a file with the name libXXX.a (or libpackage1.a).

In the above, the dot in -L. is the usual Unix pathname that denotes the current working directory. And it is important that there be no whitespace after a -L or a -l option and its value.

This syntax generalizes to multiple libraries in multiple directories as follows. Suppose that the libraries libaaa.a, libbbb.a and libccc.a are in the directory L1 and that the libraries libddd.a, libeee.a and libfff.a are in the directory L2. In this case, the link list would look like (split here into two lines):

```
-L<path-to-L1> -laaa -lbbb -lccc
-L<path-to-L2> -lddd -leee -lfff
```

The -L -l syntax is in common use throughout many Unix build systems: if your link list contains many object libraries from a single directory then it is not necessary to repeatedly specify the path to the directory; once is enough. If you are writing link lists by hand, this is very convenient. In a script, if the path name of the directory is very long, this convention makes a much more readable link list.

To perform the exercise using build3, stay in the same directory and cleanup then rebuild as follows:

1. Remove files built by build2:

```
rm function.o t1.o libpackage1.a t1
```

2. Build the code with build2 and look at the directory contents:

```
build3
ls
build build3 function.h libpackage1.a t1.cc
build2 function.cc function.o t1
```

## 3. Run ./t1 as before

The difference between build2 and build3 is that build3 compiles the main program and links it, all one one line. build2, on the other hand did the two steps separately.

## 6.6 Classes

#### 6.6.1 Introduction

The comments in the sample program used in Section 6.3 emphasized that every variable has a type: int, float, std::string, std::vector<std::string>, and so on. One of the basic building blocks of C++ is that users may define their own types; user-defined types may be built-up from all types, including other user-defined types.

The most common user-defined type is called a  $class(\gamma)$ . As you work through the Workbook exercises, you will see classes that are defined by the Workbook itself; you will also see classes defined by the toyExperiment UPS product; you will see classes defined by art and you will see classes defined by the many UPS products that support art. You will also write some classes of your own. When you work with the software for your experiment you will work with classes defined within your experiment's software.

In general, a class contains both a *declaration* (what it consists of) and an *instantiation*( $\gamma$ ) (what to do with the parts). The declaration contains some data, called *data members* or *member data*, plus some functions (called *member functions*) that will operate on that data, but it is legal for a class declaration (and therefore, a class) to contain only data or only functions. A class *declaration* has the following form:

## class MyClassName {

}

The string *class* is an identifier that is reserved to C++ and may not be used for any user-defined *identifiers*. This construct tells the C++ compiler that MyClassName is the name of a class; everything that is between the braces is part of the class declaration.

```
// required: declarations of all members of the class
// optional: definitions of some members of the class
```

The remainder of Section 6.6 will give many examples of *members* of a class.

In a class declaration, the semi-colon after the closing brace is important.



The upcoming sections will illustrate some features of classes, with an emphasis on features that will be important in the earlier Workbook exercises. This is not indended to be a comprehensive description of classes. To illustrate, we will show nine versions of a class named Point that represents a point in a plane. The first version will be simple and each subsequent version will add features.

This documentation will use technically correct language so that you will find it easier to read the standard reference materials.

#### 6.6.2 C++ Exercise 4 v1: The Most Basic Version

Here you will see a very basic version of the class Point and an illustration of how Point can be used. The ideas of *data members*, *objects* and *instantiation* will be defined.

To build and run this example:

- 1. Log in and follow the follow the steps in Section 6.2.
- 2. cd to the directory for this exercise and examine it:

```
cd Classes/v1/
ls
Point.h ptest.cc
```

Within the subdirectory v1 the main program for this exercise is the file ptest.cc. The file Point.h contains the first version of the class Point; shown in Listing ??.

3. Build the exercise.

```
../build
ls
Point.h ptest ptest.cc
```

The file named ptest is the executable program.

4. Run the exercise:

```
./ptest
p0: (2.31827e-317, 0)
p0: (1, 2)
p1: (3, 4)
p2: (1, 2)
Address of p0: 0x7fff883fe680
Address of p1: 0x7fff883fe670
Address of p2: 0x7fff883fe660
```

The values printed out in the first line of the output may be different when you run the program (remember initializaion?). When you look at the code you will see that p0 is not properly initialized and therefore contains stale data. The last three lines of output should also differ when you run the program; they are memory addresses.

Look at the header file Point. h which shows the basic version of the class Point. The three lines starting with # make up a code guard, described in Section 32.8.

```
#ifndef Point_h
                            Code guards, described in Sec-
                            tion 32.8.
#define Point_h
                         The class declaration says that the
    class Point {
                         name of the class is Point.
    public:
                      The body of the class declaration (the lines between the
      double x:
                      braces {...}) declares two data members of the class,
      double y;
                      named x and y, both of which are of type double.
                      The line public: says that the member data x and y
};
                      are accessible by any code. Instead of public, mem-
                      bers may be declared private or protected; these
                      ideas will be discussed later.
 #iendif /* Point_h *
                             A code guard
```

The plural of *data member* is sometimes written *data members* and sometimes as *member data*.

In this exercise there is no file Point.cc because the class Point consists only of a declaration; there is no implementation to put in a corresponding .cc file.

Look at the function main() (the *main program*) in ptest.cc, below, which illustrates the use of the class PointThis file includes Point.h so that the compiler will know about the class Point when it begins execution. It also includes the C++ header <iostream> which enables printing with std::cout.

```
#include "Point.h"
#include <iostream>
int main()
                Point p0; declares p0 the name of a variable whose type is (the class) Point
{
                then prints out the values of the two data members. In C++, the dot (period) char-
                acter used this way is called the member selection operator.
 Point p0;
std::cout « "p0: (" « p0.x « ", " « p0.y « ")" « std::endl;
                These lines show how to modify the values of the data members of the object
p0.x = 1.0;
                p0, then the program makes a printout to verify that the values have indeed
p0.y = 2.0;
                changed.
std::cout « "p0: (" « p0.x « ", " « p0.y « ")" « std::endl;
 Point p1;
                   These lines declare another object, named p1, of type Point and
                   assign values to its data members. These are followed by a print
p1.x = 3.0;
                   statement.
p1.y = 4.0;
std::cout « "p1: (" « p1.x « ", " « p1.y « ")" « std::endl;
                    This declares that the object named p2 is of type Point and it assigns the value
                    of p2 to be a copy of the value of p0. When the compiler sees this line, it knows
 Point p2 = p0;
                     to copy all of the data members of the class; this is a tremendous convenience for
                    classes with many data members. Again, a print statement follows.
std::cout « "p2: (" « p2.x « ", " « p2.y « ")" « std::endl;
The last three lines print the address of each of the three objects, p0, p1 and p2 in hexidecimal fomat.
std::cout « "Address of p0: " « &p0 « std::endl;
std::cout « "Address of p1: " « &p1 « std::endl;
std::cout « "Address of p2: " « &p2 « std::endl;
return 0;
 }
}
```

When the first line of code in the main () program,

### Point p0;

is executed, the program will ensure that memory has been allocated<sup>†</sup> to hold the data members of p0. If the class Point contained code to initialize data members then the program would also run that, but Point does not have any such code. Therefore the data members take on whatever values happened to preexist in the memory that was allocated for them.

Some other standard pieces of C++ nomenclature can now be defined:

- 1. The identifier p0 refers to a variable in the source code whose type is Point.
- 2. When the running program executes this line of code, it  $instantiates(\gamma)$  the  $object(\gamma)$  with the identifier p0.
- 3. The object with the identifier p0 is an instance( $\gamma$ ) of the class Point.
- 4. The identifier p0 now also refers to a region of memory containing the bytes belonging to an object of type Point.

An important take-away from the above is that a *variable* is an identifier in a source code file while an *object* is something that exists in the computer memory. Most of the time a one-to-one correspondence exists betweeen variables in the source code and objects in memory. There are exceptions, however, for example, sometimes a compiler needs to make anonymous temporary objects that do not correspond to any variable in the source code, and sometimes two or more variables in the source code can refer to the same object in memory.

The last section of the main program (and of ptest.cc itself) prints the address of each of the three objects, p0, p1 and p2. The addresses are represented in hexadecimal (base 16) format. On almost all computers, the length of a double is eight bytes. Therefore an object of type Point will have a length of 16 bytes. If you look at the printout made by ptest you will see that the addresses of p0, p01 and p2 are separated by 16 bytes; therefore the three objects are contiguous in memory.

Figure 6.1 shows a diagram of the computer memory at the end of running ptest; the outer box (blue outline) represents the memory of the computer; each filled colored box

<sup>†</sup> This is deliberately vague — there are many ways to allocate memory, and sometimes the memory allocation is actually done much earlier on, perhaps at link time or at load time.

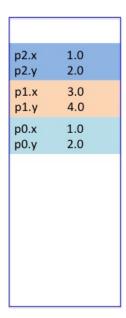


Figure 6.1: Memory diagram at the end of a run of Classes/v1/ptest.cc

represents one of the three objects in this program. The diagram shows them in contiguous memory locations, which is not necessary; there could have been gaps between the memory locations in Figure 6.1.

Now, for a bit more terminology: each of the objects p0, p1 and p2 has the three attributes required of an *object*:

- 1. a *state*, given by the values of its data members
- 2. the ability to have operations performed on it: e.g., setting/reading in value of a data member, assigning value of object of a given type to another of the same type
- 3. an *identity*: a unique address in memory

### 6.6.3 C++ Exercise 4 v2: The Default Constructor

This exercise expands the class Point by adding a default  $constructor(\gamma)$ .

To build and run this example:

Part

- 1. Log in and follow the follow the steps in Section 6.2.

```
2. Go to the directory for this exercise:
   cd Classes/v2
   ls
  Point.cc Point.h ptest.cc
  In this example, Point.cc is a new file.
3. Build the exercise:
   ../build
   ls
  Point.cc Point.h ptest ptest.cc
4. Run the exercise:
   ptest
  p0: (0, 0)
```

When you run the code, all of the printout should match the above printout exactly.

Look at Point.h. There is one new line in the body of the class declaration:

p0: (3.1, 2.7)

```
Point();
```

The parentheses tell you that this new member is some sort of function. A C++ class may have several different kinds of functions.

A function that has the same name as the class itself has a special role and is called a *con*structor; if a constructor takes no arguments it is called a default constructor. In informal written material, the word constructor is sometimes written as *c'tor*.



Point.h declares that the class Point has a default constructor, but does not define

it (i.e., provide an implementation). The definition/implementation of the constructor is found in the file Point.cc.

Look at the file Point.cc. It "includes" the header file Point.h because the compiler needs to know all about this class before it can compile the code that it finds in Point.cc. The rest of the file contains a *definition* of the constructor. The syntax Point:: says that the function to the right of the:: is part of (a member of) the class Point. The body of the constructor gives initial values to the two data members, x and y.

Look at the program ptest.cc. The first line of the main program is again

```
Point p0;
```

When the program executes this line, the first step is the same as before: it ensures that memory has been allocated for the data members of p0. This time, however, it also calls the default constructor of the class Point (declared in Point.h), which initializes the two data members (per Point.cc) such that they have well defined initial values. This is reflected in the printout made by the next line.

The next block of the program assigns new values to the data members of p0 and prints them out.

In the previous example, Classes/v1/ptest.cc, a few things happened behind the scenes that will make more sense now that you know what a constructor is.

- 1. Since the class Point did not contain a default constructor, the compiler (implicitly) wrote a default constructor for you; this default constructor simply "default constructed" each of the data members.
- 2. The (implicit) constructor of the built-in type double did nothing, leaving the data members x and y uninitialized.

### 6.6.4 C++ Exercise 4 v3: Constructors with Arguments

This exercise introduces four new ideas:

- 1. constructors with arguments
- 2. the copy constructor

- 3. implicitly generated constructor
- 4. single-phase construction vs. two-phase construction

To build and run this exercise, cd to the directory Classes/v3 and follow the same instructions as in Section 6.6.3. When you run the ptest program, you should see the following output:

### ptest

```
p0: (1, 2) p1: (1, 2)
```

Look at the file Point.h. This contains one new line:

```
Point ( double ax, double ay);
```

This line declares a second constructor; we know it is a constructor because it is a function whose name is the same as the name of the class. It is distinguishable from the default constructor because its argument list is different than that of the default constructor. As before, the file Point.h contains only the declaration of this constructor, not its *definition* (aka *implementation*).

Look at the file Point.cc. The new content in this file is the implementation of the new constructor; it assigns the values of its arguments to the data members. The names of the arguments, ax and ay, have no meaning to the compiler; they are just identifiers. It is good practice to choose names that bear an obvious relationship to those of the data members. One convention that is sometimes used is to make the name of the argument be the same as that of the data member, but with a prefix letter a, for argument. Whatever convention you (or your experiment) choose(s), use it consistently. When you update code that was initially written by someone else, follow whatever convention they adopted. Choices of style should be made to reinforce the information present in the code, not to fight it.

Look at the file ptest.cc. The first line of the main program is now:

```
Point p0(1.,2.);
```

This line declares the variable p0 and initializes it by calling the new constructor defined in this section. The next line prints the value of the data members.

The next line of code

```
Point p1(p0);
```

introduces the *copy constructor*. A copy constructor is indicated by code (like the above) that wants to create an exact copy (e.g., p1, data-member-for-data-member, of an existing type/class (e.g., p0). This exercise does not provide a copy constructor so the compiler implicitly declares a copy constructor, with public access, for that class. (The compiler puts the constructor code directly into the object file; it does not affect the source file.)

In general<sup>‡</sup> if no user-defined constructor exists for a class A, the compiler implicitly declares a default, parameterless constructor A:A() when it needs to create an object of type A. The constructor will have no constructor initializer and a null body.

We recommend that for any class whose data members are either built-in types or simple aggregates of built-in types, of which Point is an example, you let the compiler write the copy constructor for you.

If your class has data members that are pointers, or data members that manage some external resource, such as a file that you are writing to, then you will very likely need to write your own copy constructor. There are some other cases in which you should write your own copy constructor, but discussing them here is beyond the scope of this document. When you need to write your own copy constructor, you can learn how to do so from any standard C++ reference; see Section 6.7.

The next line in the file prints the values of the data members of p1 and you can see that the copy constructor worked as expected.

Notice that in the previous version of ptest.cc, the variable p0 was initialized in three lines:

```
Point p0;
p0.x = 3.1;
p0.y = 2.7;
```

This is called *two-phase construction*. In contrast, the present version uses *single-phase construction* in which the variable p0 is initialized in one line:

```
Point p0(1.,2.);
```

<sup>‡</sup>Some text in this section is adapted from material in publib.boulder.ibm.com/infocenter.

We strongly recommend using single-phase construction whenever possible. Obviously it takes less real estate, but more importantly:



- 1. Single-phase construction more clearly conveys the intent of the programmer: the intent is to initialize the object p0. The second version says this directly. In the first version you needed to do some extra work to recognize that the three lines quoted above formed a logical unit distinct from the remainder of the program. This is not difficult for this simple class, but it can become so with even a little additional complexity.
- 2. Two-phase construction is less robust. It leaves open the possibility that a future maintainer of the code might not recognize all of the follow-on steps that are part of construction and will use the object before it is fully constructed. This can lead to difficult-to-diagnose run-time errors.

### 6.6.5 C++ Exercise 4 v4: Colon Initializer Syntax

This version of the class Point introduces *colon initializer syntax* for constructors.

To build and run this exercise, cd to the directory Classes/v4 and follow the same instructions as in the previous two sections. When you run the ptest program you should see the following output:

### ptest

```
p0: (1, 2) p1: (1, 2)
```

The file Point.h is unchanged between this version and the previous one.

Now look at the file Point.cc, which contains the *definitions* of both constructors. The first thing to look at is the default constructor, which has been rewritten using colon initializer syntax. The rules for the colon-initializer syntax are:

- 1. A colon must immediately follow the closing parenthesis of the argument list.
- 2. There must be a comma-separated list of data members, each one initialized by calling one of its constructors.

art Documentation

- 3. In the initializer list, the data members must be listed in the order in which they appear in the class declaration.
- 4. The body of the constructor, enclosed in braces, must follow the initializer list.
- 5. If a data member is missing from the initializer list, its default constructor will be called (constructors for the missing data members will be called in the order in which data members were specified in the class declaration).
- 6. If no initializer list is present, the compiler will call the default constructor of every data member and it will do so in the order in which data members were specified in the class declaration.

If you think about these rules carefully, you will see that in Classes/v3/Point.cc:

- 1. the compiler did not find an initializer list, so it wrote one that default-constructed  $\mathbf x$  and  $\mathbf y$
- 2. it then wrote the code to make the assignments x=0 and y=0

On the other hand, when the compiler compiled the code for the default constructor in Classes/v4/Point.cc, it wrote the code to construct x and y, both set to zero.

Therefore, the machine code for the v3 version does more work than that for the v4 version. In practice Point is a sufficiently simple class that the compiler likely recognized and elided all of the unnecessary steps in v3; it is likely that the compiler actually produced identical code for the two versions of the class. For a more complex class, however, the compiler may not be able to recognize meaningless extra work and it will write the machine code to do that extra work.

In many cases it does not matter which of these two ways you use to write a constructor; but on those occasions that it does matter, the right answer is always the colon-initializer syntax. So we strongly recommend that you always use the colon initializer syntax. In the Workbook, all classes are written with colon-initializer syntax.

Now look at the second constructor in Point.cc; it also uses colon-initializer syntax but it is laid out differently. The difference in layout has no meaning to the compiler — whitespace is whitespace. Choose which ever seems natural to you.

Look at ptest.cc. It is the same as the version v3 and it makes the same printout.

### 6.6.6 C++ Exercise 4 v5: Member functions

This section will introduce  $member functions(\gamma)$ , both  $const member functions(\gamma)$  and non-const member functions. It will also introduce the header <cmath>. Suggested homework for this material follows.

To build and run this exercise, cd to the directory Classes/v5 and follow the same instructions as in Section 6.6.3. When you run the ptest program you should see the following output:

### ptest

```
Before p0: (1, 2) Magnitude: 2.23607 Phi: 1.10715
After p0: (3, 6) Magnitude: 6.7082 Phi: 1.10715
```

Look at the file Point.h. Compared to version v4, this version contains three additional lines:

```
double mag() const;
double phi() const;
void scale( double factor );
```

All three lines declare *member functions*. As the name suggests, a *member function* is a function that can be called and it is a member of the class. Contrast this with a *data member*, such as x or y, which are not functions. A member function may access any or all of the member data of the class.

The member function named mag does not take any arguments and it returns a double; you will see that the value of the double is the magnitude of the 2-vector from the origin to (x,y). The identifier const represents a contract between the definition/implementation of mag and any code that uses mag; it "promises" that the implementation of mag will not modify the value of any data members. The consequences of breaking the contract are illustrated in the homework at the end of this subsection.

Similarly, the member function named phi takes no arguments, returns a double and has the const identifier. You will see that the value of the double is the azimuthal angle of the vector from the origin to the point (x, y).

The third member function, scale, takes one argument, factor. Its return type is void, which means that it returns nothing. You will see that this member function multiplies both

x and y by factor (i.e., changing their values). This function declaration does not have the const identifier because it actually does modify member data.



If a member function does not modify any data members, you should always declare it const simply as a matter of course. Any negative consequences of not doing so might only become apparent later, at which point a lot of tedious editing will be required to make everything right.

Look at Point.cc. Near the top of the file an additional include directive has been added; <math> is a header from the C++ standard library that declares a set of functions for computing common mathematical operations and transformations. Functions from this library are in the  $namespace(\gamma)$  std.

Later on in Point.cc you will find the definition of mag, which computes the magnitude of the 2-vector from the origin to (x,y). To do so, it uses std::sqrt, a function declared in the <cmath> header that takes the square root of its argument. The identifier const that was present in the declaration of mag must also be present in its definition.

The next part of Point.cc contains the definition of the member function phi. To do its work, this member function uses the atan2 function from the standard library.

The next part of Point.cc contains the definition of the member function scale. You can see that this member function simply multiplies the two data members by the value of the argument.

The file ptest.cc contains a main()) program that illustrates these new features. The first line of this function declares and initializes an object, p0, of type Point. It then prints out the value of its data members, the value returned from calling the function mag and the value returned from calling phi. This shows how to access a member function: you write the name of the variable, followed by a dot (the *member selection operator*), followed by the name of the member function and its argument list.

The next line calls the member function scale with the argument 3. The printout verifies that the call to scale had the intended effect.

One final comment is in order. Many other modern computer languages have ideas very similar to C++ classes and C++ member functions; in some of those languages, the name *method* is the technical term corresponding to *member function* in C++. The name *method* 

is not part of the formal definition of C++, but is commonly used nonetheless. In this documentation, the two terms can be considered synonymous.

Here we suggest four activities as homework to help illustrate the meaning of const and to familiarize you with the error messages produced by the C++ compiler. Before moving to a subsequent activity, undo the changes that you made in the current activity.

1. In the definition of the member function Point::mag(), found in Point.cc, before taking the square root, multiply the member datum x by 2.

```
double Point::mag() const{
  x *= 2.;
  return std::sqrt( x*x + y*y );
}
```

Then build the code again; you should see the following diagnostic message:

```
Point.cc: In member function 'double Point::mag() const':
Point.cc:13:8: error: assignment of member 'Point::x' in
    read-only object
```

2. In ptest.cc, change the first line to

```
Point const p0(1,2);
```

Then build the code again; you should see the following diagnostic message:

```
ptest.cc: In function 'int main()':
ptest.cc:13:14: error: no matching function for call to
    'Point::scale(double) const'
ptest.cc:13:14: note: candidate is:
In file included from ptest.cc:1:0:
Point.h:13:8: note: void Point::scale(double) <near match>
Point.h:13:8: note: no known conversion for implicit
    'this' parameter from 'const Point*' to 'Point*'
```

These first two homework exercises illustrate how the compiler enforces the contract defined by the identifier const that is present at the end of the declaration of Point::mag() and that is absent in the definition of the member function Point::scale(). The contract says that the definition of Point::mag() may not modify the values of any data members of the class Point; users of the class Point may count on this behaviour.

The contract also says that the definition of the member function Point::scale() may modify the values of data members of the class Point; users of the class Point must assume that Point::scale() will indeed modify member data and act accordingly.§

In the first homework exercise, the value of a member datum is modified, thereby breaking the contract. The compiler detects it and issues a diagnostic message.

In the second homework exercise, the variable p0 is declared const; therefore the code may not call non-const member functions of p0, only const member functions. When the compiler sees the call to p0.mag() it recognizes that this is a call to const member function and compiles the call; when it sees the call to p0.scale(3.) it recognizes that this is a call to a non-const member function and issues a diagnostic message.

4. In Point.h, remove the const identifier from the declaration of the member function Point::mag():

```
double mag();
```

Then build the code again; you should see the following diagnostic message:

5. In Point.cc, remove the const identifier in definition of the member function mag. Then build the code again; you should see the following diagnostic message:

The third and fourth homework exercises illustrate that the compiler considers two member functions that are identical except for the presence of the const identifier to be different

<sup>§</sup> C++ has another identifier, mutable, that one can use to exempt individual data members from this contract. It's use is beyond the scope of this introduction and it will be described when it is encountered.

functions \( \). In homework exercise 3, when the compiler tried to compile Point::mag() const in Point.cc, it looked at the class declaration in Point.h and could not find a matching member function declaration; these was a close, but not exact match. Therefore it issued a diagnostic message, telling us about the close match, and then stopped. Similarly, in homework exercise 4, it also could not find a match.

### 6.6.7 C++ Exercise 4 v6: Private Data and Accessor Methods

### 6.6.7.1 Setters and Getters

This version of the class Point is used to illustrate the following ideas:

- 1. The class Point has been redesigned to have private data members with access to them provided by *accessor functions* and *setter functions*.
- 2. the *this pointer*
- 3. Even if there are many objects of type Point in memory, there is only one copy of the code.

A 2D point class, with member data in Cartesian coordinates, is not a good example of *why* it is often a good idea to have private data. But it does have enough richness to illustrate the mechanics, which is the purpose of this section. Section 6.6.7.3 discusses an example in which having private data makes obvious sense.

To build and run this exercise, cd to the directory Classes/v6 and follow the same instructions as in Section 6.6.3. When you run the ptest program you should see the following output: ptest

```
Before p0: (1, 2) Magnitude: 2.23607 Phi: 1.10715
After p0: (3, 6) Magnitude: 6.7082 Phi: 1.10715
p1: (0, 1) Magnitude: 1 Phi: 1.5708
p1: (1, 0) Magnitude: 1 Phi: 0
p1: (3, 6) Magnitude: 6.7082 Phi: 1.10715
```

Look at Point.h. Compare it to the version in v5:

Another way of saying the same thing is that the const identifier is part of the  $signature(\gamma)$  of a function.

### diff -wb Point.h ../v5/

Relative to version v5 the following changes were made:

1. four new member functions have been declared.

```
(a) double x() const;
(b) double y() const;
(c) void set( double ax, double ay);
(d) void set( Point const& p);
```

- 2. the data members have been declared private
- 3. the data members have been renamed from x and y to x\_ and y\_

Yes, there are two functions named set. Since in C++ the full name of a member function encodes all of the following information:

- 1. the name of the class it is in
- 2. the name of the member function
- 3. the argument list; that is the number, type and order of arguments
- 4. whether or not the function is const.

the member functions both named set are completely different member functions. As you work through the Workbook you will encounter a lot of this and you should develop the habit of looking at the full function name (i.e., all the parts). The full name of a member function, turned into text string, is called the *mangled name* of the member function; each C++ compiler does this a little differently. All linker symbols related to C++ classes are the mangled names of the members.

If you want to see what mangled names are created for the class Point, you can do the following



```
c++ -Wall -Wextra -pedantic -Werror -std=c++11 -c Point.cc
```

nm Point.o You can understand the output of nm by reading the man page for nm.

Part

In a class declaration, if any of the identifiers public, private, or protected appear, then all members following that identifier, and before the next such identifier, have the named property. In Point. h the two data members are private and all other members are public.

Look at Point.cc. Compare it to the version in v5:

```
diff -wb Point.cc ../v5/
```

Relative to version v5 the following changes were made:

- 1. the data members have been renamed from x and y to x\_ and y\_
- 2. an implementation is present for each of the four new member functions

Inspect the code in the implementation of each of the new member functions. The member function x () simply returns the value of the data member x; similarly for the member function y (). These are called *accessors*, *accessor functions*, or *getters*  $\|$  . The notion of *accessor* is often extended to include any member function that returns the value of simple, non-modifying calculations on a subset of the member data; in this sense, the mag and phi functions of the Point class are considered *accessors*.

The two member functions named set copy the values of their arguments into the data members of the class. These are, not surprisingly, called *setters* or *setter functions*.

More generally, any member function that modifies the value of any member data is called a *modifier*.

There is no requirement that there be accessors and setters for every data member of a class; indeed, many classes provide no such member functions for many of their data members. If a data member is important for managing internal state but is of no value to a user of the class, then you should certainly not provide an accessor or a setter.

Now that the data members of Point are private, i.e., only the code within Point is permitted to access these data members directly. All other code must access this information via the accessor and setter functions.

There is a coding style in which the function x() would have been called something like GetX(), getX() or  $get_x()$ ; hence the name *getters*. Almost all of the code that you will see in the Workbook omits the get in the names of accessors; the authors of this code view the get as redundant. Within the Workbook, the exception is for accessors defined by ROOT. The Geant4 package also includes the Get in the names of its accessors.

```
Point::Point(): x_(0.), y_(0.){}
p1.x
          3.0
                             Point::x() const { return x_;}
p1.y
          6.0
p0.x
          3.0
                             Point::y() const { return y_;}
p0.y
          6.0
                              Point::Point(double ax, double ay):
                                  x_{ax}, y_{ay}
                              double Point::mag() const {
                              return std::sqrt( x_*x_ + y_*y_);
                              double Point::phi() const {
                              return std::atan2(y_,x_);
```

**Figure 6.2:** Memory diagram at the end of a run of Classes/v6/ptest.cc

Look at ptest.cc. Compare it to the version in v5:

```
diff -wb ptest.cc ../v5/
```

Relative to version v5 the following changes were made:

- 1. the printout has been changed to use the accessor functions
- 2. a new section has been added to illustrate the use of the two set methods

Presumably these are clear.

Figure 6.2 shows a diagram of the computer memory at the end of running this version of ptest. The two boxes with the blue outlines represent sections of the computer memory; the part on the left represents that part that is reserved for storing data (such as objects) and the part on the right represents the part of the computer memory that holds the executable code. This is a big oversimplification because, in a real running program, there are many parts of the memory reserved for different sorts of data and many parts reserved for executable code.

The key point in Figure 6.2 is that each object has its own member data but there is only one copy of the code. Even if there are thousands of objects of type Point, there will only be one copy of the code. When a line of code asks for p0.mag(), the computer will pass the address of p0 as an argument to the function mag(), which will then do its work. When a line of code asks for p1.mag(), the computer will pass the address of p1 as an argument to the function mag(), which will then do its work.

Intially this sounds a little weird: the previous paragraph talks about passing an argument to the function mag() but, according to the source code, mag() does not take any arguments! The answer is that all member functions have an implied argument that always must be present — the address of the object that the member function will do work on. Because it must always be there, and because the compiler knows that it must always be there, there is no point in actually writing it in the source code! It is by using this so called hidden argument that the code for mag() knew that x means one thing for p0 but that it means something else for p1.

Every C++ member function has a variable whose name is this, which is a pointer to the object on which the member function will do its work. For example, the accessor for x () could have been written:

```
double x() const { return this->x_; }
```

This version of the syntax makes it much clearer how there can be one copy of the code even though there are many objects in memory; but it also makes the code harder to read once you have understood how the magic works. There are not many places in which you need to explicitly use the *this* pointer, but there will be some. For further information, consult standard C++ documentation (listed in Section 6.7).

### 6.6.7.2 What's the deal with the underscore?

C++ will not permit you to use the same name for both a data member and its accessor. Since the accessor is part of the public interface, it should get the simple, obvious, easy-to-type name. Therefore the name of the data member needs to be decorated to make it distinct.

The convention used in the Workbook exercises and in the toyExperiment UPS product is that the names of member data end in an underscore character. There are some other

conventions that you may encounter:

```
__name;
__name;
m_name;
mName;
theName;
```



You may also see the choice of a leading underscore, or double underscore, followed by a capital letter. Never do this.



The compiler promises that all of the linker symbols it creates will begin with a leading single or double underscore, followed by a capital letter. Some of the identifiers that you define in a C++ class will be used as part of a linker symbol. If you chose identifiers that match the pattern reserved for symbols created by the compiler there is a chance you will have naming collision with a compiler defined symbol. While this is a very small risk, it seems wise to adopt habits that guarantee that it can never happen.

It is common to extend the pattern for decorating the names of member data to all member data, even those without accessors. One reason for doing so is just symmetry. A second reason has to do with writing member functions; the body of a member function will, in general, use both member data and variables that are local to the member function. If the member data are decorated differently than the local variables, it can make the member functions easier to understand.

### 6.6.7.3 An example to motivate private data

This section describes a class for which it makes sense to have private data: a 2D point class that has data members r and phi instead of x and y. The author of such a class might wish to define a standard representation in which it is guaranteed that r be non-negative and that phi be on the domain  $0 <= \phi < 2\pi$ . If the data is public, the class cannot make these guarantees; any code can modify the data members and break the guarantee.

If this class is implemented with private data manipulated by member functions, then the constructors and member functions can enforce the guarantees.

The language used in the software engineering texts is that a guaranteed relationship among the data members is called an *invariant*. If a class has an invariant then the class

must have private data.

If a class has no invariant then one is free to choose public data. The Workbook and the toyExperiment never make this choice. One reason is that classes that begin life without an invariant sometimes acquire one as the design matures — we recommend that you plan for this unless you are 100% sure that the class will never have an invariant. A second the reason is that mixing private and public data is very confusing when you're starting out.

### 6.6.8 C++ Exercise 4 v7: The inline Identifier

This section introduces the inline identifier.

To build and run this exercise, cd to the directory Classes/v7 and follow the same instructions as in Section 6.6.3. When you run the ptest program you should see the following output: ptest

```
p0: (1, 2) Magnitude: 2.23607 Phi: 1.10715
```

Look at Point.h and compare it to the version in v6. The new material added to this version is the implementation for the two accessors x() and y(). These accessors are defined outside of the class declaration.

Look at Point.cc and compare it to the version in v6. You will see that the implementation of the accessors x() and y() has been removed.

Point.h now contains an almost exact copy of the the implementation of the accessor x () that was previously found in the file Point.cc; the difference is that it is now preceded by the identifier inline. This identifier tells the compiler that it has two options that it may choose from at its discretion.

The first option is that the compiler may decline to write a callable member function x (); instead, whenever the member function x () is used, the compiler will insert the body of x () right into the machine code at that spot. This is called *inlining* the function. For something simple like an accessor, relative to explicitly calling a function, the inlined code is very likely to

1. have a smaller memory footprint

### 2. execute more quickly

These are both good things.

On the other hand, if you inline a bigger or more complex function, some negative effects of inlining may appear. If the inlined function is used in many places and if the memory footprint of the inlined code is large compared to the memory footprint of a function call, then the total size of the program can increase. There are various ways in which a large program might run more slowly than a logically equivalent but smaller program. So, if you inline large functions, your program may actually run more slowly!

When the compiler sees the inline identifier, it also has a second option: it can choose to ignore it. When the compiler chooses this option it will write many copies of the code for the member function — one copy for each *compilation unit\*\** in which the function is called. Each compilation unit only knows about its own copy of the function and the compiler calls that copy as needed. The net result is completely negative: the function call is not actually elided so there is no time savings from that; moreover the code has become bigger because there are multiple copies of the function in memory; the larger memory footprint can further slow down execution; and compilation takes longer because multiple copies of the function must be compiled.

C++ does not permit you to force inlining; you may only give a hint to the compiler that a function is appropriate for inlining.

The bottom line is that you should always inline simple accessors and simple setters. Here the adjective *simple* means that they do not do any significant computation and that they do not contain any if statements or loops. The decision to inline anything else should only follow careful analysis of information produced by a profiling tool.

Look at the definition of the member function y() in Point.h. Compared to the definition of the member function x() there is small change in whitespace. This difference is not meaningful to the compiler. You will see several other variations on whitespace when you look at code in the Workbook and its underlying packages.

<sup>\*\*</sup> A compilation unit is the unit of code that the compiler considers at one time. For most purposes, each .cc file is its own compilation unit.

# 6.6.9 C++ Exercise 4 v8: Defining Member Functions within the Class Declaration

The version of Point in this section introduces the idea that you may provide the definition (implementation) of a member function at the point that it is declared inside the class declaration. This topic is introduced now because you will see this syntax as you work through the Workbook.

To build and run this exercise, cd to the directory Classes/v8 and follow the same instructions as in Section 6.6.3. When you run the ptest program you should see the following output:

### ptest

```
p0: (1, 2) Magnitude: 2.23607 Phi: 1.10715
```

This is the same output made by v7.

Look at Point.h. The only change relative to v7 is that the definition of the accessor methods x() and y() has been moved into the class declaration.

The files Point.cc and ptest.cc are unchanged with respect to v7.

This version of Point.h shows that you may define any member function inside the class declaration. When you do this, the inline identifier is implicit. Section 6.6.8 discussed some cautions about inappropriate use of inlining; those same cautions apply when a member function is defined inside the class declaration.

When you define a member function within the class declaration, you must not prefix the function name with the class name and the scope resolution operator; that is,



```
double Point::x() const { return x_; }
```

would produce a compiler diagnostic.

In summary, there are two ways to write inlined definitions of member functions. In most cases, the two are entirely equivalent and the choice is simply a matter of style. The one exception occurs when you are writing a class that will become part of an *art* data product, in which case it is recommended that you write the definitions of member functions *outside* of the class declaration.





When writing an *art* data product, the code inside that header file is parsed by software that determines how to write objects of that type to the output disk files and how to read objects of that type from input disk files. The software that does the parsing has some limitations and we need to work around them. The work arounds are easiest if any member functions definitions in the header file are placed outside of the class declarations. For details see

https://cdcvs.fnal.gov/redmine/projects/art/wiki/Data\_Product\_Design\_Guide #Issues-mostly-related-to-ROOT

### 6.6.10 C++ Exercise 4 v9: The stream insertion operator

The version of Point in this section illustrates how to write a *stream insertion operator*. This is the piece of code that lets you print an object without having to print each data member by hand, for example:

```
Point p0(1,2);
std::cout << p0 << std::endl;</pre>
```

To build and run this exercise, cd to the directory Classes/v9 and follow the same instructions as in Section 6.6.3. When you run the ptest program you should see the following output:

### ptest

```
p0: (1, 2) Magnitude: 2.23607 Phi: 1.10715
```

This is the same output made by v7 and v8.

Look at Point.h. The changes relative to v7 are the following two additions:

- 1. an include directive for the header <iosfwd>
- 2. a declaration for the stream insertion operator

Look at Point.cc. The changes relative to v7 are the following two additions:

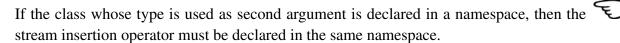
- 1. an include directive for the header <iostream>
- 2. the definition of the stream insertion operator.

Part

Look at ptest.cc. The only change relative to v7 is that the printout now uses the stream insertion operator for p0 instead of inserting each data member of p0 by hand.

In Point.h, the stream insertion operator is declared as (shown here on two lines)

```
std::ostream& operator<<
(std::ostream& ost, Point const& p );</pre>
```



When the compiler sees a « operator that has an object of type std::ostream on its left hand side and an object of type Point on its right hand side, then the compiler will look for a function named operator« whose first argument is of type std::ostream& and whose second argument is of type Point const&. If it finds such a function it will call that function to do the work; if it cannot find such a function it will issue a compiler diagnostic.

The reason that the function returns a std::ostream& is that this is the C++ convention that permits us to chain together multiple instances of the « operator:

```
Point p0(1,2), p1(3,4);
std::cout << p0 << '' '' << p1 << std::endl;
```

The C++ compiler parses this left to right. First it recognizes:

```
std::cout << p0;
```

and calls our stream insertion operator to do this work. Then it thinks of the rest of the line as:

```
std::cout << '' '' << p1 << std::endl;
Now it recognizes,</pre>
```

```
std::cout << '' ';
```

and calls the appropriate stream insertion operator to do that work. And so on.

Look at the implementation of the stream insertion operator in Point.cc. The first argument, ost, is a reference to an object of type output stream; the name ost has no meaning to C++; it is just a variable. When writing this operator we don't know and don't care what

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the output stream is connected to; perhaps it is a file; perhaps it is standard output. In any case, you send output to ost just as you do to std::cout, which is just another object of type std::ostream. In this example we chose to enclose the values of  $x_a$  and  $y_a$  in parentheses and to separate them with a comma; this is simply our choice, not something required by C++ or by art.

In this example, the stream insertion operator does *not* end by inserting a newline into ost. This is a very common choice as it allows the user of the operator to have full control about line breaks. For a class whose printout is very long and covers many lines, you might decide that this operator should end by inserting newline character; it's your choice.

If you wish to write a stream insertion operator for another class, just follow the pattern used here.

If you want to understand more about why the operator is written the way that it is, consult the standard C++ references; see Section 6.7.

The stream insertion operator is a *free function*( $\gamma$ ), not a member function of the class Point; the tie to the class Point is via its second argument. Because this function is a free function, it could have been declared in its own header file and its implementation could be provided in its own .cc file. However that is not common practice. Instead the common practice is as shown in this example: to include it in Point.h and Point.cc.



The choice of whether or not to put the declaration of the stream insertion operator into its own header file is a tradeoff between the following two criteria:

- 1. it is convenient to have it there; otherwise you would have to remember to include an additional header file when you want to use this operator
- 2. one can imagine many simple free functions that take an object of type Point as an argument. If we put them all inside Point.h, and if they are only infrequently used, then the compiler will waste time processing those declarations every time Point.h is included somewhere.

Ultimately this is a judgement call and the code in this example follows the recommendations made by the *art* development team. Their recommendation is that the following sorts of free functions, and only these sorts, should be included in header files containing a class

### declaration:

- 1. the stream insertion operator for that class
- 2. out of class arithmetic and comparison operators

With one exception, if including a function declaration in Point. h requires the inclusion of an additional header in Point.h, declare that function in a different header file. The exception is that it is okay to include <iosfwd>.

### 6.6.11 Review

The class Point is an example of a class that is primarily concerned with providing convenient access to the data it contains. Not all classes are like this; when you work through the Workbook, you will write some classes that are primarily concerned with packaging convenient access to a set of related functions.

- 1. class
- 2. object
- 3. identifier
- 4. free function
- 5. member function

## 6.7 C++ References

This section lists some recommended C++ references, both text books and online materials.

The following references describe the C++ core language,

- Stroustrup, Bjarne: "The C++ Programming Language, Special Third Edition", Addison-Wesley, 2000. ISBN 0-201-70073-5.
- http://www.cplusplus.com/doc/tutorial/

The following references describe the C++ Standard Library,

- Josuttis, Nicolai M., "The C++ Standard Library: Tutorial and Reference", Addison-Wesley, 1999. ISBN 0-201-37926-0.
- http://www.cplusplus.com/reference

The following contains an introductory tutorial. Many copies of this book are available at the Fermilab library. It is a very good introduction to the big ideas of C++ and Object Oriented Programming but it is not a fast entry point to the C++ skills needed for HEP.

 Andrew Koenig and Barbara E. Moo, "Accelerated C++: Practical Programming by Example" Addison-Wesley, 2000. ISBN 0-201-70353-X.

The following contains a discussion of recommended best practices,

Herb Sutter and Andrei Alexandrescu, "C++ Coding Standards: 101 Rules, Guidelines, and Best Practices.", Addison-Wesley, 2005. ISBN 0-321-11358-6.

## 7 Using External Products in UPS

Section 3.6.8 introduced the idea of external products. For the Intensity Frontier experiments (and for Fermilab-based experiments in general), access to external products is provided by a Fermilab-developed product-management package called Unix Product Support (UPS). An important UPS feature – demanded by most experiments as their code evolves – is its support for multiple versions of a product and multiple builds (e.g., for different platforms) per version.

Another notable feature is its capacity to handle multiple databases of products. So, for example, on Fermilab computers, login scripts (see Section 4.9) set up the UPS system, providing access to a database of products commonly used at Fermilab.



The *art* Workbook and your experiment's code will require additional products (available in other databases). For example, each experiment will provide a copy of the toyExperiment product in its experiment-specific UPS database.

In this chapter you will learn how to see which products UPS makes available, how UPS handles variants of a given product, how you use UPS to initialize a product provided in one of its databases and about the environment variables that UPS defines.

### 7.1 The UPS Database List: PRODUCTS

The act of setting up UPS defines a number of environment variables (discussed in Section 7.5), one of which is PRODUCTS. This particularly important environment variable merits its own section.

The environment variable PRODUCTS is a colon-delimited list of directory names, i.e., it is a path (see Section 4.6). Each directory in PRODUCTS is the name of a *UPS database*,

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meaning simply that each directory functions as a repository of information about one or more products. When UPS looks for a product, it checks each directory in PRODUCTS, in the order listed, and takes the first match.



If you are on a Fermilab machine, you can look at the value of PRODUCTS just after logging in, before sourcing your site-specific setup script. Run printenv:

### printenv PRODUCTS

It should have a value of

```
/grid/fermiapp/products/common/db
```

This generic Fermilab UPS database contains a handful of software products commonly used at Fermilab; most of these products are used by all of the Intensity Frontier Experiments. This database does not contain any of the experiment-specific software nor does it contain products such as  $ROOT(\gamma)$ ,  $Geant4(\gamma)$ , CLHEP or art. While these last few products are indeed used by multiple experiments, they are often custom-built for each experiment and as such are distributed via the experiment-specific (i.e., separate) UPS databases.

After you source your site-specific setup script, look at PRODUCTS again. It will probably contain multiple directories, thus making many more products available in your "site" environment. For example, on the DS50+Fermilab site, after running the DS50 setup script, PRODUCTS contains:

```
/ds50/app/products/:grid/fermiapp/products/common/db
```

You can see which products PRODUCTS contains by running Is on its directories, one-by-one, e.g.,

### Is /grid/fermiapp/products/common/db

```
afs git ifdhc mu2e python ... cpn gitflow jobsub_tools oracle_tnsnames ... encp gits login perl setpath ...
```

### Is /ds50/app/products

```
art cetpkgsupport g4neutronxs libxml2 ... artdaq clhep g4nucleonxs messagefacility ...
```

Part

```
art_suitecmakeg4photonmpich...art_workbook_basecpp0xg4piimvapich2...boostcppunitg4radiativepython...caencommds50daqg4surfaceroot...
```

Each directory name in these listings corresponds to the name of a UPS product. If you are on a different experiment, the precise contents of your experiment's product directory may be slightly different. Among other things, both databases contain a subdirectory named ups\*; this is for the UPS system itself. In this sense, all these products, including *art*, toyExperiment and even the product(s) containing your experiment's code, regard UPS as just another external product.

## 7.2 UPS Handling of Variants of a Product

An important feature of UPS is its capacity to make multiple variants of a product available to users. This of course includes different versions, but beyond that, a given version of a product may be built more than one way, e.g., for use by different operating systems (what UPS distinguishes as *flavors*). For example, a product might be built once for use with SLF5 and again for use with SLF6. A product may be built with different versions of the C++ compiler, e.g., with the production version and with a version under test. A product may be built with full compiler optimization or with the maximum debugging features enabled. Many variants can exist. UPS provides a way to select a particular build via an idea named *qualifiers*.

The full identifier of a UPS product includes its product name, its version, its flavor and its full set of qualifiers. In Section 7.3, you will see how to fully identify a product when you set it up.

## 7.3 The setup Command: Syntax and Function

Any given UPS database contains several to many, many products. To select a product and make it available for use, you use the setup command.

<sup>\*</sup>ups appears in both listings; as always, the first match wins!

In most cases the correct flavor can be automatically detected by setup and need not be specified. However, if needed, flavor, in addition to various qualifiers and options can be specified. These are listed in the UPS documentation referenced later in this section. The version, if specified, must directly follow the product name in the command line, e.g.,:

```
setup <options> <product-name> <product-version> -f <flavor> -q <qualifiers>
```

Putting in real-looking values, it would look something like:

```
setup -R myproduct v3_2 -f SLF5 -q BUILD_A
```

What does the setup command actually do? It may do any or all of the following:

- o define some environment variables
- o define some bash functions
- o define some aliases
- o add elements to your PATH
- o setup additional products on which it depends

Setting up dependent products works recursively. In this way, a single setup command may trigger the setup of, say, 15 or 20 products.

When you follow a given site-specific setup procedure, the PRODUCTS environment variable will be extended to include your experiment-specific UPS repository.



setup is a bash function (defined by the UPS product when it was initialized) that shadows a Unix system-configuration command also named setup, usually found in /usr/bin/setup or /usr/sbin/setup. Running the right 'setup' should work automatically as long as UPS is properly initialized. If it's not, setup returns the error message:

You are attempting to run ''setup'' which requires administrative privileges, but more information is needed in order to do so.

If this happens, the simplest solution is to log out and log in again.

Few people will need to know more than the above about the UPS system. Those who do can consult the full UPS documentation at:

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http://www.fnal.gov/docs/products/ups/ReferenceManual/index.html

### 7.4 Current Versions of Products

For some UPS products, but not all, the site administrator may define a particular fully-qualified version of the product as the default version. In the language of UPS this notion of default is called the *current* version. If a current version has been defined for a product, you can set up that product with the command:

### setup product-name>

When you run this, the UPS system will automatically insert the version and qualifiers of the version that has been declared current.

Having a current version is a handy feature for products that add convenience features to your interactive environment; as improvements are added, you automatically get them.

However the notion of a current version is very dangerous if you want to ensure that software built at one site will build in exactly the same way on all other sites. For this reason, the Workbook fully specifies the version number and qualifiers of all products that it requires; and in turn, the products used by the Workbook make fully qualified requests for the products on which they depend.



## 7.5 Environment Variables Defined by UPS

When your login script or site-specific setup script initializes UPS, it defines many environment variables in addition to PRODUCTS (Section 7.1), one of which is UPS\_DIR, the root directory of the currently selected version of UPS. The script also adds \$UPS\_DIR/bin to your PATH, which makes some UPS-related commands visible to your shell. Finally, it defines the bash function setup (see Sections 4.8 and 7.3). When you use the setup command, as illustrated below, it is this bash function that does the work.

In discussing the other important variables, the toyExperiment product will be used as an example product. For a different product, you would replace "toyExperiment" or "TOY-EXPERIMENT" in the following text by the product's name. Once you have followed your appropriate setup procedure (Table 5.1) you can issue the following command this

is informational for the purposes of this section; you don't need to do it until you start running the first Workbook exercise):

### setup toyExperiment v0\_00\_15 -qe2:prof

The version and qualifiers shown here are the ones to use for the Workbook exercises. When the setup command returns, the following environment variables will be defined:

TOYEXPERIMENT\_DIR defines the root DIRectory of the chosen UPS product

TOYEXPERIMENT\_INC defines the path to the root directory of the C++ header files that are provided by this product (so called because the header files are INCluded)

TOYEXPERIMENT\_LIB defines the directory that contains all of the shareable object LIBraries (ending in .so) that are provided by this product

Almost all UPS products that you will use in the Workbook define these three environment variables. Several, including toyExperiment, define many more. Once you're running the exercises, you will be able to see all of the environment variables defined by the toyExperiment product by issuing the following command:

### printenv | grep TOYEXPERIMENT



Many software products have version numbers that contain dot characters. UPS requires that version numbers not contain any dot characters; by convention, version dots are replaced with underscores. Therefore v0.00.14 becomes v0\_00\_14. Also by convention, the environment variables are all upper case, regardless of the case used in the product names.

## 7.6 Finding Header Files

### 7.6.1 Introduction

*Header files* were introduced in Section 6.3.2. Recall that a header file typically contains the "parts list" for its associated .cc source file and is "included" in the .cc file.

The software for the Workbook depends on a large number of external products; the same is true, on an even larger scale, for the software in your experiment. The preceding sec-

tions in this chapter discussed how to establish a working environment in which all of these software products are available for use.

When you are working with the code in the Workbook, and when you are working on your experiment, you will frequently encounter C++ classes and functions that come from these external products. An important skill is to be able to identify them when you see them and to be able to follow the clues back to their source and documentation. This section will describe how to do that.

An important aid to finding documentation is the use of *namespaces*; if you are not familiar with namespaces, see Section 32.6 or consult the standard C++ documentation.

### 7.6.2 Finding art Header Files

This subsection will use the example of the class art::Event to illustrate how to find header files from the *art* UPS product; this will serve as a model for finding header files from most other UPS products.

The class that holds the *art* abstraction of an HEP event is named, art::Event; that is, the class Event is in the namespace art. In fact, all classes and functions defined by *art* are in the namespace art. The primary reason for this is to minimize the chances of accidental name collisions between *art* and other codes; but it also serves a very useful documentation role and is one of the clues you can use to find header files.

If you look at code that uses art:: Event you will almost always find that the file includes the following header file:

```
#include "art/Framework/Principal/Event.h"
```

The *art* UPS product has been designed so that the relative path used to include any *art* header file starts with the directory art; this is another clue that the class or function of interest is part of *art*.

When you setup the *art* UPS product, it defines the environment variable ART\_INC, which points to the root of the header file tree for *art*. You now have enough information to discover where to find the header file for art::Event; it is at

```
$ART_INC/art/Framework/Principal/Event.h
```

You can follow this same pattern for any class or function that is part of *art*. This will only work if you are in an environment in which ART\_INC has been defined, which will be described in Chapters 9 and 10.

If you are new to C++, you will likely find this header file difficult to understand; you do not need to understand it when you first encounter it but, for future reference, you do need to know where to find it.

Earlier in this section, you read that if a C++ file uses art::Event, it would *almost always* include the appropriate header file. Why *almost* always? Because the header file Event.h might already be included within one of the other headers that are included in your file. If Event.h is indirectly included in this way, it does not hurt also to include it explicitly, but it is not required that you do so.†

We can summarize this discussion as follows: if a C++ source file uses art::Event it must always include the appropriate header file, either directly or indirectly.

art does not rigorously follow the pattern that the name of file is the same as the name of the class or function that it defines. The reason is that some files define multiple classes or functions; in most such cases the file is named after the most important class that it defines

Finally, from time to time, you will need to dig through several layers of header files to find the information you need.

There are two code browsing tools that you can use to help navigate the layering of header files and to help find class declarations that are not in a file named for the class:

- 1. use the *art redmine*( $\gamma$ ) repository browser: https://cdcvs.fnal.gov/redmine/projects/art/repository/revisions/master/show/art
- 2. use the LXR code browser: http://cdcvs.fnal.gov/lxr/art/

(In the above, both URLs are live links.)

<sup>†</sup> Actually there is small price to pay for redundant includes; it makes the compiler do unnecessary work, and therefore slows it down. But providing some redundant includes as a pedagodical tool is often a good trade-off; the Workbook will frequently do this.

### 7.6.3 Finding Headers from Other UPS Products

Section 3.7 introduced the idea that the Workbook is built around a UPS product named toyExperiment, which describes a made-up experiment. All classes and functions defined in this UPS product are defined in the namespace tex, which is an acronym-like shorthand for toyExperiment (ToyEXperiment). (This shorthand makes it (a) easier to focus on the name of each class or function rather than the namespace and (b) quicker to type.)

One of the classes from the toyExperiment UPS product is tex::GenParticle, which describes particles created by the event generator, the first part of the simulation chain (see Section 3.7.2). The include directive for this class looks like

```
#include "toyExperiment/MCDataProducts/GenParticle.h"
```

As for headers included from *art*, the first element in the relative path to the included file is the name of the UPS product in which it is found. Similarly to *art*, the header file can be found using the environment variable TOYEXPERIMENT\_INC:

```
$TOYEXPERIMENT INC/toyExperiment/MCDataProducts/GenParticle.h
```

With a few exceptions, discussed in Section 7.6.4, if a class or function from a UPS product is used in the Workbook code, it will obey the following pattern:

- 1. The class will be in a namespace that is unique to the UPS product; the name of the namespace may be the full product name or a shortened version of it.
- 2. The lead element of the path specified in the include directive will be the name of the UPS product.
- 3. The UPS product setup command will define an environment variable named <PRODUCT-NAME>\_INC, where <PRODUCT-NAME> is in all capital letters.

Using this information, the name of the header file will always be

```
$<PRODUCT-NAME>_INC/<path-specified-in-the-include-directive>
```

This pattern holds for all of the UPS products listed in Table 7.1.

A table listing git- and LXR-based code browsers for many of these UPS products can be found near the top of the web page:

https://cdcvs.fnal.gov/redmine/projects/art/wiki

**Table 7.1:** For selected UPS Products, this table gives the names of the associated namespaces. The UPS products that do not use namespaces are discussed in Section 7.6.4. <sup>‡</sup>The namespace tex is also used by the *art* Workbook, which is not a UPS product.

UPS Product	Namespace
art	art
boost	boost
cet	cetlib
clhep	CLHEP
fhiclepp	fhicl
messagefacility	mf
toyExperiment	tex <sup>‡</sup>

### 7.6.4 Exceptions: The Workbook, ROOT and Geant4

There are three exceptions to the pattern described in Section 7.6.3:

- o the Workbook itself
- o ROOT
- o Geant4

The Workbook is so tightly coupled to the toyExperiment UPS product that all classes in the Workbook are also in its namespace, tex. Note, however, that classes from the Workbook and the toyExperiment UPS product can still be distinguished by the leading element of the relative path found in the include directives for their header files:

- o art-workbook for the Workbook
- toyExperiment for the toyExperiment

The ROOT package is a CERN-supplied software package that is used by *art* to write data to disk files and to read it from disk files. It also provides many data analysis and data presentation tools that are widely used by the HEP community. Major design decisions for ROOT were frozen before namespaces were a stable part of the C++ language, therefore ROOT does not use namespaces. Instead ROOT adopts the following conventions:

1. All class names by defined by ROOT start with the capital letter T followed by another upper case letter; for example, TFile, TH1F, and TCanvas.

- 2. With very few exceptions, all header files defined by ROOT also start with the same pattern; for example, TFile.h, TH1F.h, and TCanvas.h.
- 3. The names of all global objects defined by ROOT start with a lower case letter g followed by an upper case letter; for example gDirectory, gPad and gFile.

The rule for writing an include directive for a header file from ROOT is to write its name without any leading path elements:

```
#include "TFile.h"
```

All of the ROOT header files are found in the directory that is pointed to by the environment variable \$ROOT\_INC. For example, to see the contents of this file you could enter:

```
less $ROOT_INC/TFile.h
```

Or you can the learn about this class using the reference manual at the CERN web site: http://root.cern.ch/root/html534/ClassIndex.html

You will not see the Geant 4 package in the Workbook but it will be used by the software for your experiment, so it is described here for completeness. Geant 4 is a toolkit for modeling the propagation particles in electromagnetic fields and for modeling the interactions of particles with matter; it is the core of all detector simulation codes in HEP and is also widely used in both the Medical Imaging community and the Particle Astrophysics community.

As with ROOT, Geant4 was designed before namespaces were a stable part of the C++ language. Therefore Geant4 adopted the following conventions.

- 1. The names of all identifiers begin with G4; for example, G4Step and G4Track.
- 2. All header file names defined by Geant4 begin with G4; for example, G4Step.h and G4Track.h.

Most of the header files defined by Geant4 are found in a single directory, which is pointed to by the environment variable G4INCLUDE.

The rule for writing an include directive for a header file from Geant4 is to write its name without any leading path elements:

```
#include "G4Step.h"
```

The workbook does not set up a version of Geant4; therefore G4INCLUDE is not defined. If it were, you would look at this file by:

### less \$G4INCLUDE/G4Step.h



Both ROOT and Geant4 define many thousands of classes, functions and global variables. In order to avoid collisions with these identifiers, do not define any identifiers that begin with any of (case-sensitive):

- o T, followed by an upper case letter
- o g, followed by an upper case letter
- o G4

# Part II

# Workbook

# 8 Preparation for Running the Workbook Exercises

# 8.1 Introduction

You will run the Workbook exercises on a computer that is maintained by your experiment, either at Fermilab or at another institution. Many details of the working environment change from site to site\* and these differences are parameterized so that (a) it is easy to establish the required environment, and (b) the Workbook exercises work the same way at all sites. In this chapter you will learn how to find and log into the right machine remotely from your local machine (laptop or desktop), and make sure it can support your Workbook work.



Note that is possible to install the Workbook software on your local (Unix-like) machine; instructions are available at . The instructions in this document will work whether the Workbook code is installed locally or on a remote machine.

# 8.2 Getting Computer Accounts on Workbook-enabled Machines

In order to run the exercises in the Workbook, you will need an account on a machine that can access your site's installation of the Workbook code. The experiments provide instructions for getting computer accounts on their machines (and various other information for new users) on web pages that they maintain, as listed in Table 8.1. The URLs in the table are live hyperlinks.

<sup>\*</sup>Remember, a *site* refers to a unique combination of experiment and institution.

Table 8.1: Experiment-specific Information for New Users

Experiment	URL of New User Page
ArgoNeut	https://cdcvs.fnal.gov/redmine/projects/larsoftsvn/wiki/Using_LArSoft_on_the_GPVM_nodes
Darkside	https://cdcvs.fnal.gov/redmine/projects/darkside-public/wiki/Before_You_Arrive
LArSoft	https://cdcvs.fnal.gov/redmine/projects/larsoftsvn
LBNE	https://cdcvs.fnal.gov/redmine/projects/larsoftsvn/wiki/Using_LArSoft_on_the_GPVM_nodes
MicroBoone	https://cdcvs.fnal.gov/redmine/projects/larsoftsvn/wiki/Using_LArSoft_on_the_GPVM_nodes
Muon g-2	https://cdcvs.fnal.gov/redmine/projects/g-2/wiki/NewGm2Person
Mu2e	http://mu2e.fnal.gov/atwork/general/userinfo/index.shtml#comp
NOvA	http://www-nova.fnal.gov/NOvA_Collaboration_Information/index.html

Currently, each of the experiments using *art* has installed the Workbook code on one of its experiment machines in the Fermilab General Purpose Computing Farm (GPCF).



At time of writing, the new-user instructions for all LArSoft-based experiments are at the LArSoft site; there are no separate instructions for each experiment.

If you would like a computer account on a Fermilab computer in order to evaluate *art*, contact the *art* team (see Section 3.4).

# 8.3 Choosing a Machine and Logging In

The experiment-specific machines confirmed to host the Workbook code are listed in Table 8.2 In most cases the name given is not the name of an actual computer, but rather a round-robin alias for a cluster. For example, if you log into mu2evm, you will actually be connected to one of the five computers mu2egpvm01 through mu2egpvm05. These Mu2e machines share all disks that are relevant to the Workbook exercises, so if you need to log in multiple times, it is perfectly OK if you are logged into two different machines; you will still see all of the same files.

Each experiment's web page has instructions on how to log in to its computers from your local machine.

art Documentation

Experiment	Name of Login Node
ArgoNeut	argoneutvm.fnal.gov
Darkside	ds50.fnal.gov
LBNE	lbnevm.fnal.gov
MicroBoone	uboonevm.fnal.gov
Muon g-2	gm2gpvm.fnal.gov
Mu2e	mu2evm.fnal.gov
$NO\nu A$	nova-offline.fnal.gov

**Table 8.2:** Login machines for running the Workbook exercises

# 8.4 Launching new Windows: Verify X Connectivity

Some of the Workbook exercises will launch an X window from the remote machine that opens in your local machine. To test that this works, type xterm &:

### xterm &



This should, without any messages, give you a new command prompt. After a few seconds, a new shell window should appear on your laptop screen; if you are logging into a Fermilab computer from a remote site, this may take up to 10 seconds. If the window does not appear, or if the command issues an error message, contact a computing expert on your experiment.

To close the new window, type exit at the command prompt in the new window:

# exit



If you have a problem with xterm, it could be a problem with your Kerberos and/or ssh configurations. Try logging in again with ssh -Y.

# 8.5 Choose an Editor

As you work through the Workbook exericses you will need to edit files. Familiarize yourself with one of the editors available on the computer that is hosting the Workbook. Most Fermilab computers offer four reasonable choices: emacs, vi, vim and nedit. Of these, nedit is probably the most intuitive and user-friendly. All are very powerful once you have learned to use them. Most other sites offer at least the first three choices. You can always

contact your local system administrator to suggest that other editors be installed.

A future version of this documentation suite will include recommended configurations for each editor and will provide links to documentation for each editor.

# 9 Exercise 1: Run Pre-built art Modules

# 9.1 Introduction

In this first exercise of the Workbook, you will be introduced to the  $FHiCL(\gamma)$  configuration language and you will run *art* on several modules that are distributed as part of the toyExperiment UPS product. You will not compile or link any code.

# 9.2 Prerequisites

Before running any of the exercises in this Workbook, you need to be familiar enough with the material discussed in Part I (Introduction) of this documentation set and Chapter 8 to be able to find information as needed.

If you are following the instructions below on a Mac computer, and if you are reading the instructions from a PDF file, be aware that if you use the mouse or trackpad to cut and paste text from the PDF file into your terminal window, the underscore characters will be turned into spaces. You will have to fix them before the commands will work.



# 9.3 What You Will Learn

In this exercise you will learn:

- o when to use the site-specific setup procedure
- o a little bit about the art run-time environment (Section 9.4)
- how to set up the toyExperiment UPS product (Section 9.6.1)

Part

- how to run an *art* job (Section 9.6.1)
- how to control the number of events to process (Section ??
- how to select different input files (Section ??
- o how to start at a run, subRun or event that is not the first one in the file (Section 9.8.6)
- how to concatenate input files (Section ??
- o how to write an output file (Section ??
- o some basics about the grammar and structure of a FHiCL file (Section 9.8)
- o how art finds modules and configuration (FHiCL) files. (Sections 9.10 and 9.11)

# 9.4 The art Run-time Environment

This discussion is aimed to help you understand the process described in this chapter as a whole and how the pieces fit together in the *art* run-time environment. This environment is summarized in Figure 9.1. In this figure the boxes refer either to locations in memory or to files on a disk.

At the center of the figure is a box labelled "art executable;" this represents the art main program resident in memory after being loaded. When the art executable starts up, it reads its run-time configuration (FHiCL) file, represented by the box to its left. Following instructions from the configuration file, art will load shared libraries from toyExperiment, from art, from ROOT, from CLHEP and from other UPS products. All of these shared libraries (.so files) will be found in the appropriate UPS products in LD\_LIBRARY\_PATH, which points to directories in the UPS products area (box at upper right). Also following instructions from the FHiCL file, art will look for input files (box labeled "Event-data input files" at right). The FHiCL file will tell art to write its event-data and histogram output files to a particular directory (box at lower right).

One remaining box in the figure (at right, second from bottom) is not encountered in the first Workbook exercise but has been provided for completeness. In most *art* jobs it is necessary to access experiment-related geometry and conditions information; in a mature experiment, these are usually stored in a database that stands apart from the other elements in the picture.

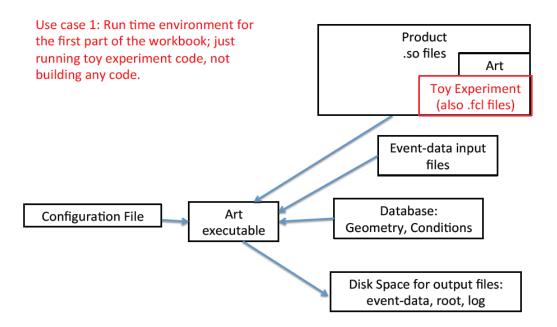


Figure 9.1: Elements of the art run-time environment for the first Workbook exercise

The arrows in Figure 9.1 show the direction in which information flows. Everything but the output flows into the *art* executable.

# 9.5 The Input and Configuration Files for the Workbook Exercises

Several event-data input files have been provided for use by the Workbook exercises. These input files are packaged as part of the toyExperiment UPS product. Table 9.1 lists the range of event IDs found in each file. You will need to refer back to this table as you proceed.

A run-time configuration (FHiCL) file has been provided for each exercise. For Exercise 1 it is hello.fcl.

Part

File Name SubRun Range of Event Numbers Run input01 data.root 1 0  $1 \dots 10$ input02 data.root 0 1...10input03\_data.root 0  $1 \dots 5$  $1 \dots 5$ 1 3 2  $1 \dots 5$ input04\_data.root 4 0  $1 \dots 1000$ 

**Table 9.1:** The input files provided for the Workbook exercises

# 9.6 Setting up to Run Exercise 1

# 9.6.1 Log In and Set Up

The intent of this section is for the reader to start from "zero" and execute an *art* job, without necessarily understanding each step, just to get familiar with the process. A detailed discussion of what these steps do will follow in Section 9.9.

Some steps are written as statements, others as commands to issue at the prompt. Notice that *art* takes the argument -c hello.fcl; this points *art* to the run-time configuration file that will tell it what to do and where to find the "pieces" on which to operate.

Most readers: Follow the steps in Section 9.6.1.1, then proceed directly to Section 9.7.

If you wish to manage your working directory yourself, skip Section 9.6.1.1, follow the steps in Section 9.6.1.2, then proceed to Section 9.7.



If you log out and wish to log back in to continue this exercise, follow the procedure outlined in Section 10.5.

# 9.6.1.1 Initial Setup Procedure using Standard Directory

- 1. Log in to the computer you chose in Section 8.3.
- 2. Follow the site-specific setup procedure; see Table 5.1.

3. Make the standard working directory then cd to it; substitute your kerberos principal for the string <username>:

```
\label{local_workbook_workbook_workbook} \begin{subarray}{l} mkdir -p $ART\_WORKBOOK\_WORKING\_BASE/<username>/ \\ \begin{subarray}{l} workbook-tutorial/pre-built \end{subarray}
```

```
{\tt cd \$ART\_WORKBOOK\_WORKING\_BASE/<username>/ \setminus workbook-tutorial/pre-built}
```

4. Setup the toyExperiment UPS product:

```
setup toyExperiment v0_00_15 -q$ART_WORKBOOK_QUAL:prof
```

5. Copy the scripts into your working directory:

```
cp $TOYEXPERIMENT_DIR/HelloWorldScripts/* .
```

6. Use the provided script to create the symbolic links needed by the FHiCL files:

```
source makeLinks.sh
```

7. See what you have in the directory:

```
ls
```

```
helloExample.log inputFiles makeLinks.sh
skipEvents.fcl
hello.fcl inputs.txt output writeFile.fcl
```

Proceed to Section 9.7.



# 9.6.1.2 Initial Setup Procedure allowing Self-managed Working Directory

- 1. Log in to the computer you chose in Section 8.3.
- 2. Follow the site-specific setup procedure; see Table 5.1
- 3. Make a working directory and cd to it.
- 4. Setup the toyExperiment UPS product:

```
setup toyExperiment v0_00_15 -q$ART_WORKBOOK_QUAL:prof
```

5. Copy the scripts into your working directory:

```
cp $TOYEXPERIMENT_DIR/HelloWorldScripts/* .
```

- 6. Make a subdirectory named output. If you prefer, you can make this on some other disk and put a symbolic link to it from the current working directory; name the link output.
- 7. Create a symbolic link to allow the FHiCL files to find the input files:

```
In -s $TOYEXPERIMENT_DIR/inputFiles .
```

8. See what you have in the directory:

```
ls
```

```
helloExample.log inputFiles makeLinks.sh
skipEvents.fcl
hello.fcl inputs.txt output writeFile.fcl
```

Proceed to Section 9.7.

### 9.6.1.3 Setup for Subsequent Exercise 1 Login Sessions

If you log out and later wish to log in again to work on this exercise, you need to do the following:

- 1. Log in to the computer you chose in Section 8.3.
- 2. Follow the site-specific setup procedure; see Section 5.
- 3. cd to your working directory, e.g., for the standard case:

```
{\tt cd \$ART\_WORKBOOK\_WORKING\_BASE/<username>/ \setminus workbook-tutorial/pre-built} \\
```

4. Setup the toyExperiment UPS product:

```
setup toyExperiment v0_00_15 -q$ART_WORKBOOK_QUAL:prof
```

Compare this with the list given in Section 9.6.1. You will see that three steps are missing because they only need to be done the first time.

You are now ready to run *art* as you were before.

# 9.7 Execute art and Examine Output

From your working directory, execute *art* on the FHiCL file hello.fcl and send the output to output/hello.log:

```
art -c hello.fcl >& output/hello.log
```

Compare the outtut you produced against Listing 9.1; the only differences should be the timestamps. *art* will have processed the first file listed in Table 9.1.

**Listing 9.1:** Sample output from running hello.fcl

```
1
2 %MSG-i MF_INIT_OK: art 27-Apr-2013 21:22:13 CDT JobSetup
3 Messagelogger initialization complete.
```

```
4 %MSG
5 27-Apr-2013 21:22:14 CDT Initiating request to open file
6 inputFiles/input01_data.root
7 27-Apr-2013 21:22:14 CDT Successfully opened file
8 inputFiles/input01_data.root
9 Begin processing the 1st record. run: 1 subRun: 0 event: 1 at
10 27-Apr-2013 21:22:14 CDT
11 Hello World! This event has the id: run: 1 subRun: 0 event: 1
12 Begin processing the 2nd record. run: 1 subRun: 0 event: 2 at
13 27-Apr-2013 21:22:14 CDT
14 Hello World! This event has the id: run: 1 subRun: 0 event: 2
15 Hello World! This event has the id: run: 1 subRun: 0 event: 3
16 Hello World! This event has the id: run: 1 subRun: 0 event: 4
17 Hello World! This event has the id: run: 1 subRun: 0 event: 5
18 Hello World! This event has the id: run: 1 subRun: 0 event: 6
19 Hello World! This event has the id: run: 1 subRun: 0 event: 7
20 Hello World! This event has the id: run: 1 subRun: 0 event: 8
21 Hello World! This event has the id: run: 1 subRun: 0 event: 9
22 Hello World! This event has the id: run: 1 subRun: 0 event: 10
23 27-Apr-2013 21:22:14 CDT Closed file inputFiles/input01_data.root
24
25 TrigReport ----- Event Summary -----
26 TrigReport Events total = 10 passed = 10 failed = 0
27
28 TrigReport ----- Modules in End-Path: e1 -----
29 TrigReport Trig Bit# Visited Passed Failed
                                                            Error Name
                                       10
30 TrigReport
               0 0
                               10
                                                                 0 hi
31
32 TimeReport ----- Time Summary ---[sec]----
33 TimeReport CPU = 0.004000 Real = 0.002411
34
35 Art has completed and will exit with status 0.
```

Every time you run *art*, the first thing to check is the last line in your output or log file. It should be Art has completed and will exit with status 0. If the status is not 0, or if this line is missing, it is an error; please contact the *art* team as described in Section 3.4.

A future version of these instructions will specify how much disk space is needed, including space for all outtut files.

# 9.8 Understanding the Configuration

The file hello.fcl gives art its run-time configuration. This file is written in the Fermilab Hierarchical Configuration Language (FHiCL, pronounced "fickle"), a language that was developed at Fermilab to support run-time configuration for several projects, including art. By convention, files written in FHiCL end in .fcl. As you work through the Workbook, the features of FHiCL that are relevant for each exericse will be explained.

art accepts some command line options that can be used in place of items in the FHiCL file. You will encounter some of these in this section.



The full details of the FHiCL language, plus the details of how it is used by *art*, are given in the Users Guide, Chapter 25. Most people will find it much easier to follow the discussion in the Workbook documentation than to digest the full documentation up front.

# 9.8.1 Some Bookkeeping Syntax

In a FHiCL file, the start of a comment is marked by the hash sign character (#); a comment may begin in any column.

The hash sign has one other use. If the first eight characters of a line are exactly #include, followed by whitespace and a quoted list of file paths, then the line will be interpreted as an *include directive* and the line containing it will be replaced by the contents of the file named in the include directive.

The basic element of FHiCL is the definition, which has the form

```
name : value
```



A group of FHiCL definitions delimited by braces  $\{\}$  is called a *table*( $\gamma$ ). Within *art*, a FHiCL table gets turned into a C++ object called a *parameter set*( $\gamma$ ); this document set will often refer to a FHiCL table as a parameter set.

The fragment of hello.fcl shown below contains the FHiCL table that configures the  $source(\gamma)$  of events that art will read in and operate on.

```
source : {
```

}

At the outermost scope of the FHiCL file, *art* will interpret the source parameter set as the description of the source of events for this run of *art*.

# module\_type: RootInput

module\_type is an identifier that tells *art* the name of a module to load and run, RootInput in this case. RootInput, a standard source module provided by *art*, reads disk files containing event-data written in *art*-defined ROOT-based format.

```
fileNames : [
"inputfiles/input01_data.root"

1
```

The string fileNames defined in the RootInput module scope gives the module a list of filenames from which to read events.

The name source is an *identifier* in art; i.e., the name source has no special meaning to FHiCL but it does have a special meaning to art. To be precise, it only has a special meaning to art if it is at the outermost  $scope(\gamma)$  of a FHiCL file; i.e., not inside any braces  $\{\}$  within the file. The notion of scope in FHiCL is discussed further in Chapter 13. When art sees a parameter set named source at the outermost scope, then art will interpret that parameter set to be the description of the source of events for this run of art.



In the source parameter set, module\_type is an identifier in *art* that tells *art* the name of a module that it should load and run, RootInput in this case. RootInput is one of the standard source modules provided by *art* and it reads disk files containing event-data written in an *art*-defined ROOT-based format. The default behaviour of the RootInput module is to start at the first event in the first file and read to the end of the last event in the last file.\*

The string fileNames is again an identifier, but this time defined in the RootInput module, that gives the module a list of filenames from which to read events. The list is delimited by square brackets and contains a comma-separated list of filenames. This exam-

<sup>\*</sup> In the Workbook, the only source module\_type that you will see will be RootInput. Your experiment may have a source module that reads events from the live experiment and other source modules that read files written in experiment-defined formats; for example Mu2e has a source module that reads single particle events from a text file written by G4beamline.

ple shows only one filename, but the square brackets are still required. The proper FHiCL name for a comma-separated list delimited by square brackets is a  $sequence(\gamma)$ .

In most cases the filenames in the sequence must be enclosed in quotes. FHiCL, like many other languages has the following rule: if a string contains white space or any special characters, then quoting it is required, otherwise quotes are optional.

FHiCL has its own set of special characters; these include anything *except* all upper and lower case letters, the numbers 0 through 9 and the underscore character. *art* restricts the use of the underscore character in some circumstances; these will be discussed as they arise.

It is implied in the foregoing discussion that a FHiCL *value* need not be a simple thing, such as a number or a quoted string. For example, in Listing ??, the source value is a parameter set (of two parameters) and the value of fileNames is a (single-item) sequence.

# 9.8.2 Some Physics Processing Syntax

The identifier  $physics(\gamma)$ , when found at the outermost scope, is an identifier reserved to art. The physics parameter set is so named because it contains most of the information needed to describe the physics workflow of an art job.

The fragment of hello.fcl below shows a rather long-winded way of telling *art* to find a module named HelloWorld and execute it.

```
At the outermost scope of the FHiCL file, art will interpret the
physics : {
                physics parameter set as the description of the physics workflow
                for this run of art.
 analyzers : {
     hi: {
     module_type: HelloWorld
                As a top-level identifier within the physics scope, analyzers defines
 }
                for art the run-time configuration for all the analyzer modules in the job,
                e.g., HelloWorld.
 e1 : [ hi ]
 end_paths: [e1]
          el and end_paths represent paths in art. el is an arbitrary identifier for a
          FHiCL sequence of module labels; end_paths, an identifier reserved to art, is
}
          a FHiCL sequence of paths.
```

Why so long-winded? *art* has very powerful features that enable execution of multiple complex chains of modules; the price is that specifying something simple takes a lot of keystrokes.

Within the physics parameter set, notice the identifier analyzers. When found as a top-level identifier within the physics scope, it is recognized as a keyword reserved to *art*. The analyzers parameter set defines the run-time configuration for all of the analyzer modules that are part of the job – only HelloWorld in this case.

For our current purposes, the module HelloWorld does only one thing of interest, namely for every event it prints one line (shown here as three):

where RR, SS and EE are substituted with the actual run, subRun and event number of each event.

If you look back at Listing 9.1, you will see that this line appears ten times, once each for events 1 through 10 of run 1, subRun 0 (as expected, according to Table 9.1). The remainder of the listing is standard output generated by *art*.

The remainder of the lines in hello.fcl appears below. The line starting with  $process\_name(\gamma)$  tells art that this job has a name and that the name is "hello"; it has no real significance in these simple exericses. It becomes important when an art job creates new data products (described in User Guide Chapter 26) and writes them to a file; each data product will be uniquely identified by a four-part name, one part of which is the name of the process that created the data product. This imposes a constraint on process\\_name values: art joins the four parts of a data product name into a single string, with the underscore (\_) as a separator between fields; none of the parts (e.g., the process name) may contain additional underscores.





In an art event-data file, each data product is stored as a TBranch of a  $TTree(\gamma)$ ; the string containing the full name of the data product is used as the name of the TBranch. On readback, art must parse the name of the TBranch to recover the four individual pieces of the data product name. If one of the four parts internally contains an underscore, then art cannot reliably recover the four parts.

The services parameter set provides run-time configuration information for all *art* services.

```
#include "fcl/minimalMessageService.fcl"

process_name : hello

The identifier process_name tells art that the name of this job is hello.

services : {

message : @local::default_message }

The services parameter set provides the run-time configuration for all the required art services, in this case the message service. Its configuration is set in the file indicated in the #include directive.
```

For our present purposes, it is sufficient to know that the configuration for the message service is found inside the file that is included via the #include line. The message service controls the limiting and routing of debug, informational, warning and error messages generated by *art* or by user code. The message service does not control information written directly to std::cout or std::cerr.

# 9.8.3 art Command line Options

art supports some command line options. To see what they are, type the following command at the bash prompt:

### art -help

Note that some options have both a short form and a long form. This is a common convention for Unix programs; the short form is convenient for interactive use and the long form makes scripts more readable.

### 9.8.4 Maximum Number of Events to Process

By default *art* will read all events from all of the specified input files. You can set a maximum number of events in two ways, one way is from the command line:

```
art -c hello.fcl -n 5
art -c hello.fcl -nevts 4
```

Run each of these commands and observe their output.

The second way is within the FHiCL file. Start by making a copy of hello.fcl:

```
cp hello.fcl hi.fcl
```

Edit hi.fcl and add the following line anywhere in the source parameter set:

```
maxEvents : 3
```

By convention this is added after the fileNames definition but it can go anywhere inside the source parameter set because the order of parameters within a FHiCL table is not important. Run *art* again, using hi.fcl:

```
art -c hi.fcl
```

You should see output from the HelloWorld module for only the first three events.

To configure the file for *art* to process all the events, i.e., to run until *art* reaches the end of the input files, either leave off the maxEvents parameter or give it a value of -1.



If the maximum number of events is specified both on the command line and in the FHiCL file, then the command line takes precedence. Compare the outputs of the following commands:

```
art -c hi.fcl
art -c hi.fcl -n 5
art -c hi.fcl -n -1
```

#### 9.8.5 **Changing the Input Files**

For historical reasons, there are multiple ways to specify the input event-data file (or the list of input files) to an art job:

- within the FHiCL file's source parameter set
- o on the art command line via the -s option (you may specify one input file only)
- o on the art command line via the -S option (you may specify a text file that lists multiple input files)
- o on the art command line, after the last recognized option (you may specify one or more input files)



If input file fiames are procedence. If input file names are provided both in the FHiCL file and on the command line, the

Let's run a few examples.

We'll start with the -s command line option (second bullet). Run art without it (again), for comparison (or recall its output from Listing 9.1):

```
art -c hello.fcl
```

To see what you should expect given the following input file, check Table 9.1, then run:

```
art -c hello.fcl -s inputFiles/input02_data.root
```

Notice that the 10 events in this output are from run 2 subRun 0, in contrast to the previous printout which showed events from run 1. Notice also that the command line specification overrode that in the FHiCL file. The -s (lower case) command line syntax will only permit you to specify a single filename.

This time, edit the source parameter set inside the hi.fcl file (first bullet); change it to:

(Notice that you also added maxEvents: -1.) The names of the two input files could have been written on a single line but this example shows that, in FHiCL, newlines are treated simply as white space.

Check Table 9.1 to see what you should expect, then rerun *art* as follows:

```
art -c hi.fcl
```

You will see 20 lines from the HelloWorld module; you will also see messages from *art* at the open and close operations on each input file.

Back to the -s command-line option, run:

```
art -c hi.fcl -s inputFiles/input03_data.root
```

This will read only inputFiles/input03\_data.root and will ignore the two files specified in the hi.fcl. The output from the HelloWorld module will be the 15 events from the three subRuns of run 3.

There are several ways to specify multiple files at the command line. One choice is to use the -S (upper case) [—source-list] command line option (third bullet) which takes as its argument the name of a text file containing the filename(s) of the input event-data file(s). An example of such as file has been provided, inputs.txt. Look at the contents of this file:

### cat inputs.txt

```
inputFiles/input01_data.root
```

```
inputFiles/input02_data.root
inputFiles/input03_data.root
```

Now run art using inputs.txt to specify the input files:

```
art -c hi.fcl -S inputs.txt
```

You should see the HelloWorld output from the 35 events in the three files; you should also see the messages from *art* about the opening and closing of the three files.

Finally, you can list the input files at the end of the *art* command line (fourth bullet):

```
art -c hi.fcl inputFiles/input02_data.root inputFiles/input03_data.root
```

(Remember the Unix convention about a trailing backslash marking a command that continues on another line; see Chapter 2. ) In this case you should see the HelloWorld output from the 25 events in the two files.



In summary, there are three ways to specify input files from the command line; all of them override any input files specified in the FHiCL file. Do not try to use two or more of these methods on a single *art* command line; the *art* job will run without issuing any messages but the output will likely be different than you expect.

# 9.8.6 Skipping Events

The source parameter set supports a syntax to start execution at a given event number or to skip a given number of events at the start of the job. Look, for example, at the file skipEvents.fcl, which differs from hello.fcl by the addition of two lines to the source parameter set:

```
firstEvent : 5
maxEvents : 3
```

art will process events 5, 6, and 7 of run 1, subRun 0. Try it:

```
art -c skipEvents.fcl
```

An equivalent operation can be done from the command line in two different ways. Try the following two commands and compare the output:

```
art -c hello.fcl -e 5 -n 3
```

Part

# art -c hello.fcl -nskip 4 -n 3

You can also specify the intial event to process relative to a given event ID (which, recall, contains the run, subRun and event number). Edit hi.fcl and edit the source parameter set as follows:

```
source : {
  module_type : RootInput
  fileNames : [ "inputFiles/input03_data.root" ]
  firstRun : 3
  firstSubRun : 1
  firstEvent : 6
}
```

When you run this job, *art* will process events starting from run 3, subRun 2, event 1, – because there are only 5 events in subRun 1.

```
art -c hi.fcl
```

# 9.8.7 Identifying the User Code to Execute

Recall from Section 9.8.2 that the physics parameter set contains the physics content for the *art* job. Within this parameter set, *art* must be able to determine which (user code) modules to process. These must be referenced via *module labels*( $\gamma$ ), which as you will see, represent the pairing of a module name and a run-time configuration.

Look back at Listing ??, which contains the physics parameter set from hello.fcl. The analyzer parameter set, nested inside the physics parameter set, contains the definition:

```
hi : {
  module_type : HelloWorld
}
```

The identifier hi is a module label (defined by the user, not by FHiCL or *art*) whose value must be a parameter set that *art* will use to configure a module. The parameter set for a module label must contain (at least) a FHiCL definition of the form:

```
module_type : <module-name>
```

Here module\_type is an identifier reserved to *art* and <module-name> tells *art* the name of the module to load and execute. (Since it is within the analyzer parameter set, the module must be of type EDAnalyzer; i.e. the *base type* of <module-name> must be EDAnalyzer.)

Module labels are fully described in Section 25.5.

In this example *art* will look for a module named <code>HelloWorld</code>, which it will find as part of the toyExperiment UPS product. Section 9.10 describes how *art* uses <code><module-name></code> to find the shareable library that contains code for the <code>HelloWorld</code> module. A parameter set that is used to configure a module may contain additional lines; if present, the meaning of those lines is understood by the module itself; those lines have no meaning either to <code>art</code> or to FHiCL.

Now look at the FHiCL fragment below that starts with analyzers. We will use it to reinforce some of the ideas discussed in the previous paragraph.

art allows you to write a FHiCL file that uses a given module more than once. For example you may want to run an analysis twice, once with a loose mass cut on some intermediate state and once with a tight mass cut on the same intermediate state. In art you can do this by writing one module and making the cuts "run-time configurable." This idea will be developed further in Chapter 14.

```
analyzers : {
    When art processes this fragment it will look for a
    module named MyAnalysis and instantiate it twice...

loose: {
    module_type : MyAnalysis
    mass_cut : 20.
    }
        ... once using the parameter set labeled loose ...

tight: {
    module_type : MyAnalysis
    mass_cut : 15.
    }
        ... and once using the parameter set labeled tight.
}
```

The two instances of the module MyAnalysis are distinguished by the module labels tight and loose. *art* requires that module labels be unique within a FHiCL file. Module labels may contain only upper- and lower-case letters and the numerals 0 to 9.



In the FHiCL files in this exercise, all of the modules are analyzer modules. Since analyzers do not make data products, these module labels are nothing more than identifiers inside the FHiCL file. For producer modules, however, which *do* make data products, the module label becomes part of the data product identifier and as such has a real signficance. All module labels must conform to the same naming rules.

Within *art* there is no notion of reserved names or special names for module labels; however your experiment will almost certainly have established some naming conventions.

### 9.8.8 Paths

In the physics parameter set for hello.fcl there are two parameters that represent paths in *art*:

```
e1 : [ hi ]
end_paths : [ e1 ]
```

The path defined by the parameter e1 takes a value that is a FHiCL sequence of module labels. The name of a path is an arbitrary identifier that must be unique within a FHiCL file; it has no persistent significance and can be any legal FHiCL name.

Sometimes this documentation uses the word path in the sense of an  $art\ path(\gamma)$  (a sequence of module labels), other times path is used as a path in a file system and in yet other situations, it is used as a colon-delimited set of directory names. The use should be clear from the context.



The name end\_paths, in contrast to e1, is an identifier reserved to *art*. Its value must be a FHiCL sequence of paths – here it is a sequence of one path, e1. *reference the rules when available* When *art* processes the end\_paths definition it combines all of the path definitions and forms the set of unique module labels from all paths defined in the parameter set. In other words, it is legal in *art* for a module label to appear in more than one path; if it does, *art* will recognize this and will ensure that the module is executed only once per event.

If you put the name of a module label into the definition of end\_paths, *art* will issue an error and stop processing.



The paths listed in end\_paths may only contain module labels for analyzer and/or output modules; they may *not* contain module labels for producer or filter modules. The reason for this restriction will be discussed in Section .

What about the order of module labels in a path? Since analyzer and output modules may neither add new information to the event nor communicate with each other except via the event, the processing order is not important for the event. By definition, then, *art* may run analyzer and output modules in any order. In a simple *art* job with a single path, *art* will, in fact, run the modules in the order of appearance in the path, but do not write code that depends on execution order because *art* is free to change it.

It may seem that end\_paths could more simply have been defined as a set of module labels, eliminating the layer of the path altogether, but there is a reason. We will defer this discussion to Section .

If the end\_paths parameter is absent or defined as:

```
end_paths : [ ]
```

*art* will understand that this job has no analyzer modules and no filter modules to execute. It is legal to define a path as an empty FHiCL sequence.

As is standard in FHiCL, if the definition of end\_paths appears more than once, the last definition takes precendence.

# 9.8.9 Writing an Output File

The file writeFile.fcl gives an example of writing an output file. This file introduces the parameter set named outputs:

the inner parameter sets provides the configuration of one output module.

An *art* job may have zero or more output modules. The name RootOutput is the name of a standard *art* output module; it writes the events in memory to a disk file in an *art*-defined, ROOT-based format. Files written by the module RootOutput can be read by the module RootInput. The identifier output1 is just another module label that obeys

Part

the same rules discussed in Section 9.8.7. The string fileName is an identifier known to the RootOutput module; its value is the name of the output file that this instance of RootOutput will write.

There are many more optional parameters that can be used to configure an output module. For example, an output module can be configured to write out only selected events and/or to write out only a subset of the available data products. Optional parameters are described in Chapter .

Notice in writeFile.fcl that the path el has been extended to include the module label of the output module:

```
el : [ hi, output1 ]
```

This ensures that both modules get entered into the path for execution.

Finally, the source parameter set of writeFile.fcl is configured to read only events 4, 5, 6, and 7.

To run writeFile.fcl and check that it worked correctly:

```
art -c writeFile.fcl
ls -s output/writeFile_data.root
art -c hello.fcl -s output/writeFile_data.root
```

The first command will write the output file; the second will check the size of the output file and the last one will read back the output file and print the event IDs for all of the events in the file. You should see the HelloWorld printout for events 4, 5, 6 and 7.

# 9.9 Understanding the Process for Exercise 1

Section 9.6.1 contained a list of steps needed to run this exercise; this section will describe each of those steps in detail. When you understand what is done in these steps, you will understand the run-time environment in which *art* runs. As a reminder, the steps are listed again here:

- 1. Log in to the computer you chose in Section 8.3.
- 2. Follow the site-specific setup procedure; see Chapter 5
- mkdir -p \$ART\_WORKBOOK\_WORKING\_BASE/<username>/ \
  workbook-tutorial/pre-built

In the above and elsewhere as indicated, substitute your kerberos principal for the string <username>.

- cd \$ART\_WORKBOOK\_WORKING\_BASE/<username>/ \
  /workbook-tutorial/pre-built
- 5. setup toyExperiment v0\_00\_15 -q\$ART\_WORKBOOK\_QUAL:prof
- 6. cp \$TOYEXPERIMENT\_DIR/HelloWorldScripts/\*.
- 7. source makeLinks.sh
- 8. Run *art*:

art -c hello.fcl >& output/hello.log

Steps 1 and 4 should be self explanatory and will not be discussed further.



When reading this section, you do not need to run any of the commands given here; this is a commentary on commands that you have already run.

# 9.9.1 Follow the Site-Specific Setup Procedure (Details)

The site-specific startup procedure, described in Chapter 5, ensures that the UPS system is properly initialized and that the UPS database (containing all of the UPS products needed to run the Workbook exercises) is present in the PRODUCTS environment variable.

This procedure also defines two environment variables that are defined by your experiment to allow you to run the Workbook exercises on their computer(s):

ART\_WORKBOOK\_WORKING\_BASE the top-level directory in which users create their working directory for the Workbook exercises

ART\_WORKBOOK\_OUTPUT\_BASE the top-level directory in which users create their output directory for the Workbook exercises; this is used by the script makeLinks.sh

If these environment variables are not defined, ask a system admin on your experiment.

# 9.9.2 Make a Working Directory (Details)

On the Fermilab computers the home disk areas are quite small so most experiments ask that their collaborators work in some other disk space. This is common to sites in general, so we recommend working in a separate space as a best practice. The Workbook is designed to require it.

This step given as:

```
\label{local_mkdir} mkdir -p \ ART_WORKBOOK_WORKING_BASE/< username > / workbook-tutorial / \\ pre-built
```

creates a new directory to use as your working directory. It is defined relative to an environment variable described in Section 9.9.1. It only needs to be done the first time that you log in to work on Workbook exercises – once it's there, it's there!

If you follow the rest of the naming scheme, you will guarantee that you have no conflicts with other parts of the Workbook.

As discussed in Section 9.6.1.2, you may of course choose your own working directory on any disk that has adequate disk space.

# 9.9.3 Setup the toyExperiment UPS Product (Details)

This step is the main event in the eight-step process.

```
setup toyExperiment v0_00_14 -q$ART_WORKBOOK_QUAL:prof
```

This command tells UPS to find a product named toyExperiment, with the specified version and qualifiers, and to *setup* that product, as described in Section 7.3.

The required qualifiers may change from one experiment to another and even from one site to another within the same experiment. To deal with this, the site specific setup procedure defines the environment variable ART\_WORKBOOK\_QUAL, whose value is the qualifier string that is correct for that site.

The complete ups qualifier for toyExperiment has two components, separated by a colon: the string defined by ART\_WORKBOOK\_QUAL plus a qualifier describing the compiler optimization level with which the product was built, in this case "prof"; see Section 3.6.7 for information about the optimization levels.

Each version of the toyExperiment product knows that it requires a particular version and qualifier of the *art* product. In turn, *art* knows that it depends on particular versions of ROOT, CLHEP, boost and so on. When this recursive setup has completed, over 20 products will have been setup. All of these products define environment variables and about two-thirds of them add new elements to the environment variables PATH and LD\_LIBRARY\_PATH.

If you are interested, you can inspect your environment before and after doing this setup. To do this, log out and log in again. Before doing the setup, run the following commands:

Then setup toyExperiment and capture the environment afterwards (env.after). Compare the before and after files: the after files will have many, many additions to the environment. (The fragment | tr: \\n tells the bash shell to take the output of printenv and replace every occurrence of the colon character with the newline character; this makes the output much easier to read.)

# 9.9.4 Copy Files to your Current Working Directory (Details)

The step:

# cp \$TOYEXPERIMENT\_DIR/HelloWorldScripts/\* .

only needs to be done only the first time that you log in to work on the Workbook.

In this step you copied the files that you will use for the exercises into your current working directory. You should see these files:

```
hello.fcl makeLinks.sh skipEvents.fcl writeFile.fcl
```

# 9.9.5 Source makeLinks.sh (Details)

This step:

### source makeLinks.sh

only needs to be done only the first time that you log in to work on the Workbook. It created some symbolic links that *art* will use.

The FHiCL files used in the Workbook exercises look for their input files in the subdirectory inputFiles. This script made a symbolic link, named inputFiles, that points to:

```
$TOYEXPERIMENT_DIR/inputFiles
```

in which the necessary input files are found.

This script also ensures that there is an output directory that you can write into when you run the exercises and adds a symbolic link from the current working directory to this output directory. The output directory is made under the directory \$ART\_WORKBOOK\_OUTPUT\_BASE; this environment variable was set by the site-specific setup procedure and it points to disk space that will have enough room to hold the output of the exercises.

### **9.9.6** Run *art* (Details)

Issuing the command:

art Documentation

### art -c hello.fcl

runs the *art* main program, which is found in \$ART\_FQ\_DIR/bin. This directory was added to your PATH when you setup toyExperiment. You can inspect your PATH to see that this directory is indeed there.

# 9.10 How does art find Modules?

When you ran hello.fcl, how did art find the module HelloWorld?

It looked at the environment variable LD\_LIBRARY\_PATH, which is a colon-delimited set of directory names defined when you setup the toyExperiments product. We saw the value of LD\_LIBRARY\_PATH in Section 9.9.3; to see it again, type the following:

```
printenv LD_LIBRARY_PATH | tr : \\n
```

The output should look similar to that shown in Listing 9.2.

Listing 9.2: Example of the value of LD\_LIBRARY\_PATH

```
1 /ds50/app/products/tbb/v4_1_2/Linux64bit+2.6-2.12-e2-prof/lib
2 /ds50/app/products/sqlite/v3_07_16_00/Linux64bit+2.6-2.12-prof/lib
3 /ds50/app/products/libsigcpp/v2_2_10/Linux64bit+2.6-2.12-e2-prof/lib
4 /ds50/app/products/cppunit/v1_12_1/Linux64bit+2.6-2.12-e2-prof/lib
5 /ds50/app/products/clhep/v2_1_3_1/Linux64bit+2.6-2.12-e2-prof/lib
6 /ds50/app/products/python/v2_7_3/Linux64bit+2.6-2.12-qcc47/lib
   /ds50/app/products/libxml2/v2_8_0/Linux64bit+2.6-2.12-gcc47-prof/lib
8 /ds50/app/products/fftw/v3_3_2/Linux64bit+2.6-2.12-gcc47-prof/lib
9 /ds50/app/products/root/v5_34_05/Linux64bit+2.6-2.12-e2-prof/lib
10 /ds50/app/products/boost/v1_53_0/Linux64bit+2.6-2.12-e2-prof/lib
/ds50/app/products/cpp0x/v1_03_15/slf6.x86_64.e2.prof/lib
12 /ds50/app/products/cetlib/v1_03_15/slf6.x86_64.e2.prof/lib2
13 /ds50/app/products/fhiclcpp/v2_17_02/slf6.x86_64.e2.prof/lib
14 /ds50/app/products/messagefacility/v1_10_16/slf6.x86_64.e2.prof/lib
15 /ds50/app/products/art/v1_06_00/slf6.x86_64.e2.prof/lib
16 /ds50/app/products/toyExperiment/v0_00_14/slf6.x86_64.e2.prof/lib
17 /grid/fermiapp/products/common/prd/git/v1_8_0_1/Linux64bit-2/lib
```

Of course the leading element of each directory name, /ds50/app will be replaced by whatever is correct for your experiment. The last element in LD\_LIBRARY\_PATH is not relevant for running *art* and it may or may not be present on your machine, depending on details of what is done inside your site-specific setup procedure.

If you compare the names of the directories listed in LD\_LIBRARY\_PATH to the names of the directories listed in the PRODUCTS environment variable, you will see that all of these directories are part of the UPS products system. Moreover, for each product, the version, flavor and qualifiers are embedded in the directory name. In particular, both *art* and toyExperiment are found in the list.

If you is the directories in LD\_LIBRARY\_PATH you will find that each directory contains many shareable object libraries (.so files).

When art looks for a module named HelloWorld, it looks through the directories defined in

LD\_LIBRARY\_PATH and looks for a file whose name matches the pattern,

```
lib*HelloWorld_module.so
```

where the asterisk matches (zero or) any combination of characters. *art* finds that, in all of the directories, there is exactly one file that matches the pattern, and it is found in the directory (shown here on two lines):

The name of the file is:

```
libtoyExperiment_Analyzers_HelloWorld_module.so
```

If *art* had found no files that matched the pattern, it would have printed an error message and tried to shutdown as gracefully as possible. If *art* had found more than one file that matched the pattern, it would have printed a different error message and tried to shut down as gracefully as possible.

One of the important features of *art* is that, whenever it detects an error condition that is serious enough to stop execution, it always attempts to shut down as gracefully as possible. Among other things this means that it tries to properly close all output files. This feature is not so important when an error occurs at the start of a job but it ensures that, when an error occurs after hours of execution, your results up to the error are correct and available.



# 9.11 How does art find FHiCL Files?

This section will describe where *art* looks for FHiCL files. There are two cases: looking for the file specified by the command line argument -c and looking for files that have been included by a #include directive within a FHiCL file.

# 9.11.1 The -c command line argument

When you issued the command

#### art -c hello.fcl

art looked for a file named hello.fcl in the current working directory and found it. You may specify any absolute or relative path as the argument of the -c option. If art had not found hello.fcl in this directory it would have looked for it relative to the path defined by the environment variable FHICL\_FILE\_PATH. This is just another path-type environment variable, like PATH or LD\_LIBRARY\_PATH. You can inspect the value of FHICL\_FILE\_PATH by:

# printenv FHICL FILE PATH

```
.: $TOYEXPERIMENT_DIR
```

Acutally the output will show the translated value of the environment variable TOYEXPERIMENT\_DIR. The presence of the current working directory (dot) in the path is redundant when processing the command line argument but it is significant in the case discussed in the next section.



Some experiments have chosen to configure their version of the *art* main program so that it will not look for the command line argument FHiCL file in FHICL\_FILE\_PATH. It is also possible to configure *art* so that only relative paths, not absolute paths, are legal values of the -c argument. This last option can be used to help ensure that only version-controlled files are used when running production jobs. Experiments may enable or disable either of these options when their main program is built.

### 9.11.2 #include Files

Section 9.8 discussed Listing ??, which contains the fragments of hello.fcl that are related to configuring the message service. The first line in that listing is an include directive. *art* will look for the file named by the include directive relative to FHICL\_FILE\_PATH and it will find it in:

\$TOYEXPERIMENT\_DIR/fcl/minimalMessageService.fcl

This is part of the toyExperiment UPS product.

The version of *art* used in the Workbook does not consider the argument of the include directive as an absolute path or as a path relative to the current working directory; it only looks for files relative to FHICL\_FILE\_PATH. This is in contrast to the choice made when processing the -c command line option.



When building *art*, one may configure *art* to first consider the argument of the include directive as a path and to consider FHICL\_FILE\_PATH only if that fails.

Add a section called Review that looks at trigger paths, end paths, etc and works backwards

# 10 Exercise 2: Build and Run Your First Module

# 10.1 Introduction

In this exercise you will build and run a simple *art* module. Section 3.6.7 introduced the idea of a build system, a software package that compiles and links your source code to turn it into machine code that the computer can execute. In this chapter you will be introduced to the *art* development environment, which adds the following to the run-time environment (discussed in Section 9.4):

- 1. a build system
- 2. a source code repository
- 3. a working copy of the Workbook source code
- 4. a directory containing shared libraries created by the build system

In this and all subsequent Workbook exercises, you will use the build system used by the *art* development team, **cetbuildtools**. This system will require you to open two shell windows your local machine and, in each one, to log into the remote machine \*. The windows will be referred to as the *source window* and the *build window*:

- In the *source window* you will check out and edit source code.
- In the *build window* you will build and run code.

<sup>\*</sup>cetbuildtools requires what are called *out-of-source builds*; this means that the source code and the working space for the build system must be in separate directories.

Exercise 2 and all subsequent Workbook exercises will use the setup instructions found in Sections 10.4 and 10.5.



## 10.2 Prerequisites

Before running this exercise, you need to be familiar with the material in Part I (Introduction) of this documentation set and Chapter 9 from Part II (Workbook). Concepts that this chapter refers to include:

- o namespace
- #include directives
- o header file
- o class
- o base class
- o derived class
- o constructor
- o destructor
- o the C preprocessor
- o member function (aka method)
- o const vs non-const member function
- o argument list of a function
- o signature of a function
- o virtual function
- o pure virtual function
- o virtual class
- o pure virtual class
- o concrete class

- o declaration vs defintion of a class
- arguments passed by reference
- o arguments passed by const reference
- o notion of type: e.g., a class, a struct, a free function or a typedef
- how to write a C++ main program

In this chapter you will also encounter the C++ idea of *inheritance*. Understanding inheritance is not a prerequisite; it will be described as you encounter it in the Workbook exercises.

#### 10.3 What You Will Learn

In this exercise you will learn:

- o how to establish the art development environment
- how to checkout the Workbook exercises from the git source code management system
- how to use the **cetbuildtools** build system to build the code for the Workbook exercises
- o how include files are found
- o what a *link list* is
- o where the build system finds the link list
- what the art::Event is and how to access it
- o what the art::EventID is and how to access it
- o what makes a class an art module
- o where the build system puts the .so files that it makes

# 10.4 Initial Setup to Run Exercises: Standard Procedure

#### 10.4.1 "Source Window" Setup

Up through step 3 of the procedure in this section, the results should look similar to those of Exercise 1. Note that the directory name chosen here in the mkdir step is different than that chosen in the first exercise; this is to avoid file name collisions.

If you want to use a self-managed working directory, in step 3 make a directory of your choosing and cd to it rather than to the directory shown.



In your source window do the following:

- 1. Log in to the computer you chose in Section 8.3.
- 2. Follow the site-specific setup procedure; see Table 5.1.
- 3. Make a new working directory and cd to it:

 $\label{eq:mkdir-p} $ART\_WORKBOOK\_WORKING\_BASE/< username > / \setminus workbook \\$ 

 ${\tt cd~\$ART\_WORKBOOK\_WORKING\_BASE/< username > / workbook}$ 

- 4. Set up the source code management system git and use it to pull down the workbook code to the directory art-workbook, which will be referred to as your *source* directory. The output for each step is explained in Section 10.4.2.1:
  - (a) setup git
  - (b) git clone http://cdcvs.fnal.gov/projects/art-workbook

- (c) cd art-workbook
- (d) git checkout -b v0\_00\_18 v0\_00\_18
- 5. Source the script that sets up the environment properly:

source ups/setup\_deps -p \$ART\_WORKBOOK\_QUAL

The git commands are discussed in Section 10.4.2.1. The final step sources a script that defines a lot of environment variables — the same set that will be defined in the build window.

#### 10.4.2 Examine Source Window Setup

#### 10.4.2.1 About git and What it Did

Git is a source code management system<sup>†</sup> that is used to hold the source code for the Workbook exercises. A source code management system is a tool that looks after the book-keeping of the development of a code base; among many other things it keeps a complete history of all changes and allows one to get a copy of the source code as it existed at any time in the past. Because of git's many advanced features, many HEP experiments are moving to git. git is fully described in the git manual<sup>‡</sup>.

Some experiments set up git in their site-specific setup procedure; others do not. In running setup git, you have ensured that a working copy of git is in your PATH<sup>§</sup>.

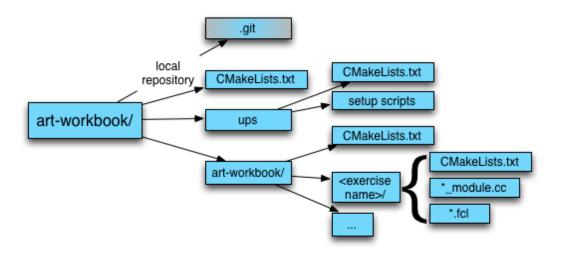
The git clone and git checkout commands produce a working copy of the Workbook source files in your source directory. Figure 10.1 shows a map of the source directory structure created by the git commands. It does not show all the contents in each subdirectory. Note that the .git (hidden) directory under the source directory is colored differently; this is done to distinguish it from the rest of the contents of the source directory structure:

<sup>†</sup>Other source code management systems with which you may be familar are CVS and SVN.

<sup>&</sup>lt;sup>‡</sup>Several references for git can be found online; the "official" documentation is found at http://git-scm.com/documentation.

<sup>&</sup>lt;sup>§</sup>No version needs to be supplied because the git UPS product has a current version declared; see Section 7.4.

- When you ran git clone in Section 10.4.1, it copied the entire contents of the remote repository into this directory. The .git directory contains your local copy of the repository.
- When you ran git checkout, it created the rest of the structure under the source directory (what we call your "working area") and copied the requested version of everything you need from .git into this structure.



**Figure 10.1:** Representation of the reader's source directory structure (an admin directory is not shown)

; git clone should produce the following output:

Cloning into 'art-workbook' ...

Executing the git checkout command should produce the following output:

Switched to a new branch 'v0\_00\_18 '

If you do not see the expected output, contact the art team as described in Section 3.4.

#### 10.4.2.2 Contents of the Source Directory

Figure 10.1 shows roughly what your source directory contains at the end of the setup procedure. You can see the correspondance between it and the output of the ls -a command:

## cd \$ART\_WORKBOOK\_WORKING\_BASE/<username>/workbook/art-workbook/s-a

```
. .. admin art-workbook CMakeLists.txt .git ups
```

Notice that it contains a subdirectory of the same name as its parent, art-workbook.

- The admin directory (not shown in Figure 10.1) contains some scripts used by **cetbuildtools** to customize the configuration. of the development environment.
- The art-workbook directory contains the main body of the source code for the Workbook exercises.
- The file CMakeLists.txt is the file that the build system reads to learn what steps it should do.
- The ups directory contains information about what UPS products this product depends on; it contains additional information used to configure the development environment.

Look inside the art-workbook (sub)directory (via ls) and see that it contains several files and subdirectories. The file CMakeLists.txt contains more instructions for the build system. Actually, every directory contains a CMakeLists.txt; each contains additional instructions for the build system. The subdirectory FirstModule contains the files that will be used in this exericse; the remaining subdirectories contain files that will be used in subsequent Workbook exercises.

If you look inside the FirstModule directory, you will see

```
CMakeLists.txt FirstAnswer01_module.cc First_module.cc firstAnswer01.fcl first.fcl
```

The file CMakeLists.txt in here contains yet more instructions for the build system and will be discussed later. The file First\_module.cc is the first module that you will look at and first.fcl is the FHiCL file that runs it. This exercise will suggest that you try to write some code on your own; the answer is provided in

FirstAnswer01\_module.cc and the file firstAnswer01.fcl runs it. These files will be discussed at length throughout the exercises.

#### 10.4.3 "Build Window" Setup

Again, advanced users wanting to manage their own working directory may skip to Section 10.4.3.2.



#### 10.4.3.1 Standard Procedure

Now go to your build window and do the following:

- 1. Log in to the computer you chose in Section 8.3.
- 2. Follow the site-specific setup procedure; see Chapter 5.
- cd \$ART\_WORKBOOK\_WORKING\_BASE/<username>/ \ workbook
- 4. mkdir build-prof

The build-prof directory will be your *build* directory.

- 5. cd build-prof
- 6. source ../art-workbook/ups/setup\_for\_development/\-p \$ART\_WORKBOOK\_QUAL

The output from this command will tell you to take some additional steps; *do not do those steps*.

7. buildtool

This step may take a few minutes.

Skip Section 10.4.3.2 and move on to Section 10.4.4.



#### 10.4.3.2 Using Self-managed Working Directory

The steps in this procedure that are the same as for the "standard" procedure are explained in Section 10.4.4.

Now go to your build window and do the following:

- 1. Log in to the computer you chose in Section 8.3.
- 2. Follow the site-specific setup procedure; see Chapter 5.
- 3. Make a directory to hold the code that you will build and cd to it; this will be your *build* directory in your *build* window.
- 4. Make another directory, *outside of the heirarchy rooted at your build directory*, to hold output files created by the workbook exercises. (Don't cd to it.)
- 5. In -s <directory-for-output-files> output
- 6. source <your-SOURCE-directory>/ups/setup\_for\_development \
  -p \$ART\_WORKBOOK\_QUAL

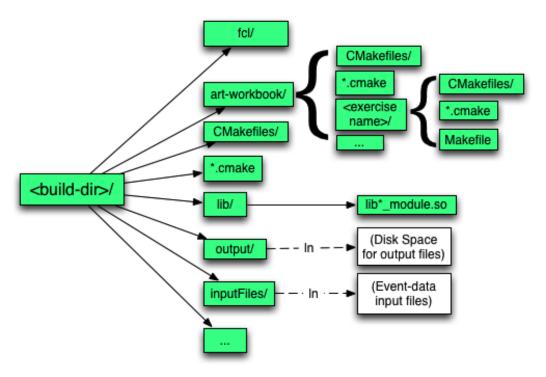
The output from this command (Listing 10.1) will tell you to take some additional steps; *do not do those steps*.

7. buildtool

#### 10.4.4 Examine Build Window Setup

Logging in and sourcing the site-specific setup script should be clear by now. Notice that next you are told to cd to the same workbook directory as in Step 4 of the instructions for the source window. From there, you make a directory in which you will run builds (your build directory), and cd to it. (The name build-prof can be any legal directory name

but it is suggested here because this example performs a profile build; this is explained in Section 3.6.7). Figure 10.2 shows roughly what the build directory contains.



**Figure 10.2:** Representation of the reader's build directory structure (the fcl/directory is a symlink to art-workbook/art-workbook/ in the source area)

Step 6 sources a script called setup\_for\_development found in the ups subdirectory of the source directory. This script, run exactly as indicated, defines build-prof to be your build directory. This command selects a profile build (via the option -p); it also requests that the UPS qualifiers defined in the environment variable ART\_WORKBOOK\_QUAL be used when requesting the UPS products on which it depends; this environment variable was discussed in Section 9.9.3. The expected output is shown in Listing 10.1.

Check that there are no error messages in the indicated block. The listing concludes with a request for you to run a cmake command; *do not run* cmake (this line is an artifact of layering **cetbuildtools** on top of cmake).



**Listing 10.1:** Example of output created by setup\_for\_development

art Documentation

```
2 The working build directory is /ds50/app/user/kutschke/workbook/build-prof
3 The source code directory is /ds50/app/user/kutschke/workbook/art-workbook
   ----- check this block for errors -----
6 /ds50/app/user/kutschke/workbook/build-prof/lib has been added to LD_LIBRARY_PATH
  /ds50/app/user/kutschke/workbook/build-prof/bin has been added to PATH
8
9 CETPKG_SOURCE=/ds50/app/user/kutschke/workbook/art-workbook
10 CETPKG BUILD=/ds50/app/user/kutschke/workbook/build-prof
11 CETPKG NAME=art workbook
12 CETPKG VERSION=v0 00 15
13 CETPKG_QUAL=e2:prof
14 CETPKG_TYPE=Prof
15
16 Please use this cmake command:
17 cmake -DCMAKE_INSTALL_PREFIX=/install/path
18
                              -DCMAKE_BUILD_TYPE=$CETPKG_TYPE $CETPKG_SOURCE
```

This script sets up all of the UPS products on which the Workbook depends; this is analogous to the actions taken by Step 6 in the first exercise (Section 9.6.1.1) when you were working in the *art* run-time environment. This script also creates several files and directories in your build-prof directory; these comprise the working space used by **cetbuild-tools**.

After sourcing this script, the contents of build-prof will be

```
art_workbook-v0_00_18 bin lib
cetpkg_variable_report diag_report
```

At this time the two subdirectories bin and lib will be empty. The other files are used by the build system to keep track of its configuration.

Step 7 (buildtool) tells **cetbuildtools** to build everything found in the source directory; this includes all of the Workbook exercises, not just the first one. The build process will take two or three mintues on an unloaded (not undergoing heavy usage) machine. Its output should end with the lines:

```
INFO: Stage build successful.
```

After the build has completed do an ls on the directory lib; you will see that it contains

Part

a large number of shared library (.so) files; for  $v0_00_18$  there will be about 30 .so files (subject to variation as versions change); these are the files that art will load as you work through the exercises.

Also do an ls on the directory bin; these are scripts that are used by **cetbuildtools** to maintain its environment; if the Workbook contained instructions to build any executable programs, they would have been written to this directory.

After runing buildtool, the build directory will contain:

admin	CMakeFiles	fcl
art-workbook	cmake_install.cmake	inputFiles
art_workbook-v0_00_15	CPackConfig.cmake	lib
bin	CPackSourceConfig.cmake	Makefile
cetpkg_variable_report	CTestTestfile.cmake	output
CMakeCache.txt	diag_report	ups

Most of these files are standard files that are explained in the **cetbuildtools** documentation. However, three of these items need special attention here because they are customized for the Workbook.

An Is -I on the files fcl, inputFiles and output will reveal that they are symbolic links to

These links are present so that the FHiCL files for the Workbook exercises can be machine-independent.

- The link inputFiles points to the directory inputFiles present in the toyExperiment UPS product; this directory contains the input files that *art* will read when you run the first exercise. These are the same files used in the first exercise; if you need a reminder of the contents of these files, see Table 9.1. These input files will also be used in many of the subsequent exercises.
- The link outputFiles points to a directory that was created to hold your output

files; the environment variable ART\_WORKBOOK\_OUTPUT\_BASE was defined by your site-specific setup procedure.

The symlink fcl points into your source directory hierarchy; it allows you to access the FHiCL files that are found in that hierarchy with the convenience of tab completions.

## 10.5 Setup for Subsequent Login Sessions

If you log out and later wish to log in again to work on this or any other subsequent exercise, you need to do the following:

In your source window:

- 1. Log in to the computer you chose in Section 8.3.
- 2. Follow the site-specific setup procedure; see Table 5.1
- 3. cd to your source directory:

```
{\tt cd \$ART\_WORKBOOK\_WORKING\_BASE/<username>/ \setminus \\ workbook/art-workbook}
```

4. Set up the environment:

```
source ups/setup_deps -p
```

In your build window:

- 1. Log in to the computer you chose in Section 8.3.
- 2. Follow the site-specific setup procedure; see Chapter 5
- 3. cd to your build directory:

```
    cd $ART_WORKBOOK_WORKING_BASE/<username>/\
workbook/build-prof
    Source the setup file:
    source ../art-workbook/ups/setup_for_development \
        -p $ART_WORKBOOK_QUAL
```

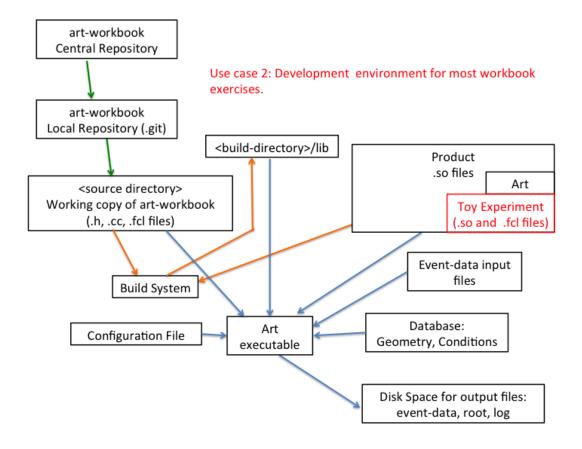
If you chose to manage your own directory names, then the names of your source and build directories will be different than those shown.

Compare these steps with those given in Sections 10.4.1 and Section 10.4.3. You will see that some steps are missing from the source window and the build window instructions. The missing steps were only required the first time.

## 10.6 The art Development Environment

In the preceding sections of this chapter you established what is known as the *art development environment*; this is a superset of the *art* run-time environment, which was described in Section 9.4. This section summarizes the new elements that are part of the development environment but not part of the run-time environment.

In Section 10.4.1, step 4b (git clone ...) contacted the central source code repository for the *art* Workbook code and made a clone of the repository in your source area under <code>art-workbook</code>; the clone contains the complete history of the repository, including all versions of art-workbook. Step 4d git checkout ...) checked out the correct version of the code from the clone and put a working copy into your source directory. The central repository is hosted on a Fermilab server and is accessed via the network. The upper left box in Figure 10.3 denotes the central repository and the box below it denotes the clone of the repository in your disk space; the box below that denotes the checked out working copy of the Workbook code. The flow of information during the clone and checkout processes is indicated by the green arrows (at left) in the figure.



**Figure 10.3:** Elements of the *art* development environment as used in most of the Workbook exercises; the arrows denote information flow, as described in the text.

In step 7 of Section 10.4.3, you ran buildtool in your build area, which read the source code files from your working copy of the Workbook code and turned them into shared libraries. The script buildtool is part of the build system, which is denoted as the box in the center left section of Figure 10.3. When you ran buildtool, it wrote shared library files to the lib subdirectory of your build directory; this directory is denoted in the figure as the box in the top center labeled <br/>
build-directory>/lib. The orange arrows in the figure denote the information flow at build-time. In order to perform this task, buildtool also needed to read header files and shared libraries found in the UPS products area, hence the orange arrow leading from the UPS Products box to the build system box.

In the figure, information flow at run-time is denoted by the blue lines. When you ran the *art* executable, it looked for shared libraries in the directories defined by LD\_LIBRARY\_PATH. In the *art* development environment, LD\_LIBRARY\_PATH contains

- 1. the lib subdirectory of your build directory
- 2. all of the directories previously described in Section 9.10

In all environments, the art executable looks for FHiCL files in

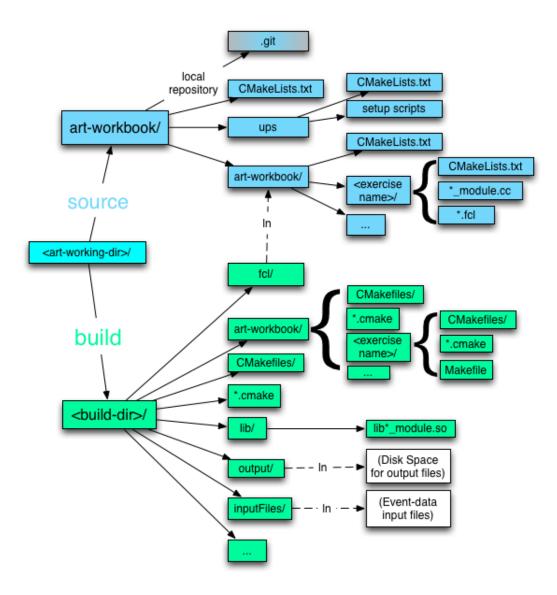
- 1. in the file specified in the -c command line argument
- 2. in the directories specified in FHICL\_FILE\_PATH

The first of these is denoted in the figure by the box labeled "Configuration File." In the art development environment, FHICL\_FILE\_PATH contains

- 1. some directories found in your checked out copy of the source
- 2. all of the directories previously described in Section 9.11

The remaining elements in Figure 10.3 are the same as described for Figure 9.1.

Figure 10.4, a combination of Figures 10.1 and 10.2), illustrates the distinct source and build areas, and the relationship between them. It does not show all the contents in each subdirectory.



**Figure 10.4:** Representation of the reader's directory structure once the development environment is established.

### 10.7 Running the Exercise

#### 10.7.1 Run art on first.fcl

In your build window, make sure that your current working directory is your build directory. From here, run the first part of this exercise by typing the following:

#### art -c fcl/FirstModule/first.fcl > output/first.log

(We suggest you get in the habit of routing your output to the output directory.) The output of this step will look much like that in Listing 9.1, but with two signficant differences. The first difference is that the output from first.fcl contains an additional line

```
Hello from First::constructor.
```

The second difference is that the words printed out for each event are a little different; the printout from first.fcl looks like

```
Hello from First::analyze. Event id: run: 1 subRun: 0 event: 1 while that from hello.fcl looked like
```

```
Hello World! This event has the id: run: 1 subRun: 0 event: 1
```

The reason for changing this printout is so that you can identify, from the printout, which module was run.

#### 10.7.2 The FHiCL File first.fcl

Compare the FHiCL file used in this exercise, fcl/FirstModule/first.fcl, with hello.fcl from the first exercise (i.e., run cat or diff on them). Other than comments, the only difference is that the module\_type has changed from HelloWorld to First:

## diff \$TOYEXPERIMENT\_DIR/HelloWorldScripts/hello.fcl \ fcl/FirstModule/first.fcl

```
composite the control of the co
```

The file first.fcl tells *art* to run a module named First. As described in Section 9.10, *art* looks through the directories defined in LD\_LIBRARY\_PATH and looks for a file whose name matches the pattern lib\*First\_module.so. This module happens to be found at this location, relative to your build directory:

```
lib/libart-workbook_FirstModule_First_module.so
```

This shared library file was created when you ran buildtool.

#### 10.7.3 The Source Code File First module.cc

This section will describe the source code for the module First and will use it as a model to describe modules in general. The source code for this module is found in the following file, relative to your source directory (go to your source window!):

```
art-workbook/FirstModule/First_module.cc
```

When you ran buildtool, it compiled and linked this source file into the following shared library (relative to your your build directory):

```
lib/libart-workbook_FirstModule_First_module.so
```

This is the shared library that was loaded by *art* when you ran code for this exercise, in Section 10.7.2.

Look at the file First\_module.cc, shown in Listing 10.2. In broad strokes, it does the following:

- o declares a class named First
- o provides the implementation for the class

Part

contains a call to the C-Preprocessor macro named DEFINE\_ART\_MODULE, discussed in Section 10.7.3.8

All module files that you will see in the Workbook share these "broad strokes." Some experiments that use *art* have chosen to split the source code for one module into three separate files; the *art* team does not recommend this practice, but it is in use and it will be discussed in Section 10.10.2.

Listing 10.2: The contents of First\_module.cc

```
1
2 #include "art/Framework/Core/EDAnalyzer.h"
3 #include "art/Framework/Core/ModuleMacros.h"
4 #include "art/Framework/Principal/Event.h"
6 #include <iostream>
8
   namespace tex {
9
10
     class First : public art::EDAnalyzer {
11
12
     public:
13
14
       explicit First(fhicl::ParameterSet const& );
15
16
       void analyze(art::Event const& event) override;
17
18
     };
19
20
21
   tex::First::First(fhicl::ParameterSet const& pset) : art::EDAnalyzer(pset) {
23
     std::cout << "Hello from First::constructor." << std::endl;</pre>
24
25
26 void tex::First::analyze(art::Event const& event) {
27
     std::cout << "Hello_from_First::analyze._Event_id:_"</pre>
28
               << event.id()
29
                << std::endl;
30 }
31
32 DEFINE_ART_MODULE(tex::First)
```

We will examine this file in the following sections.

#### 10.7.3.1 The #include Statements

```
#include "art/Framework/Core/EDAnalyzer.h"
#include "art/Framework/Core/ModuleMacros.h"
#include "art/Framework/Principal/Event.h"
```

The first three lines of code in the file First\_module.cc are *include directives* that pull in header files. All three of these files are included from the *art* UPS product (determining the location of included header files is discussed in Section 7.6).

#include <iostream>

The next line, #include <iostream>, includes the C++ header that enables this code to write output to the screen; for details, see any standard C++ documentation.

Those of you with some C++ experience may have noticed that there is no file named First\_module.h in the directory art-workbook/FirstModule. The explanation for this will be given in Section 10.10.1.



If you are a C++ beginner you will likely find these header files difficult to understand; you do not need to understand them at this time but you do need to know where to find them for future reference.

#### 10.7.3.2 The Declaration of the Class First, an Analyzer Module

Let's start with short explanations of each line and follow up with more information.

Part

```
Open a namespace named tex.
namespace tex {
 class First : public art::EDAnalyzer{
         • The first line of the declaration of the class First.
         • Analyzer modules must inherit publicly from the base class EDAnalyzer.
 public:
               Class members below here are public; any above would be private.
     explicit First(fhicl::ParameterSet const& );
         • Declaration of a constructor.
         • Its argument list is prescribed by art.
         • art will call the constructor once at the start of each job.
     void analyze(art::Event const& event) override;
         • Declaration of the analyze member function.
         • Its argument list is prescribed by art.
         • art will call this member function once per event.
         • The override contextual identifier is an important safety feature. Use it!
         Close the declaration of the class First.
 };
}
         Close the namespace tex.
```

All of the code in the toyExperiment UPS product was written in the namespace tex; the name tex is an acronym-like shorthand for the toyExperiment (ToyEXperiment) UPS product. In order to keep things simple, all of the classes in the Workbook are also declared in the namespace tex. For more information about this choice, see Section 7.6.4. If you are not familiar with namespaces, consult the standard C++ documentation.

In the first line of the class declaration, the fragment ": public art::EDAnalyzer" tells that compiler that the class First is a *dervied class* that *inherits publicly* from the *base class* art::EDAnalyzer. At this time it is not necessary to understand inheritance, which is fortunate, because it takes a long, long time to explain. You just need to recognize and follow the pattern. You can read about C++ inheritance in the standard C++

documentation.

#### 10.7.3.3 An Introduction to Analyzer Modules

Section 3.6.3 discussed the idea of *module types*: analyzer, producer, filter and so on. For a class to be a valid *art* analyzer module, it must follow a set of rules defined by *art*:

- 1. It must inherit publicly from art::EDAnalyzer.
- 2. It must provide a constructor with the argument list:

```
fhicl::ParameterSet const& pset
```

(Only the type of the argument is presecribed, not its name. You can use any name you want but the same name must be used in item 3.)

3. The initializer list of the constructor must call the constructor of the base class; and it must pass the parameter set to the constructor of the base class:

```
art::EDAnalyzer(pset)
```

4. It must provide a member function named analyze, with the signature  $\P$ :

```
analyze( art::Event const&)
```

5. If the name of a module class is <ClassName> then the source code for the module must be in a file named <ClassName>\_module.cc and this file must contain the lines:

```
#include "art/Framework/Core/ModuleMacros.h"
and
DEFINE_ART_MODULE(<namespace>::<ClassName>)
```

6. It may optionally provide other member functions with signatures prescribed by *art*; if these member functions are present in a module class, then *art* will call them at the appropriate times. Some examples are provided in Chapter 12.

A module may also contain any other member data and any other member functions that are needed to do its job. You can see from Listing 10.2 that the class First follows all of the above rules and that it does not contain any of the optional member functions.

<sup>¶</sup> In C++ the *signature* of a member function is the name of the class of which the function is a member, the name of the function, the number, types and order of the arguments, and whether the member function is marked as const or volatile. The signature does not include the return type; nor does it include the names of any of the arguments.

The requirement that the class name match the filename (minus the \_module.cc portion) is enforced by *art*'s system for dynamically loading shared libraries. The requirement that the class provide the prescribed constructor is enforced by the macro DEFINE\_ART\_MODULE, which will be described in Section 10.7.3.8.

The declaration of the constructor begins with the keyword explicit. This is a safety feature this relevant only for constructors that have exactly one argument. A proper explanation would take too long so just follow a simple guideline: all constructors that have exactly one argument should be declared explicit. There will be rare circumstances in which you need to go against this guideline but you will not encounter any in the Workbook.

The *override* contextual identifier in the analyzer member function definition is a feature that is new in C++ 11 so older references will not discuss it. It is a new safety feature that we recommend you use; we cannot give a proper explanation until we have had a chance to discuss inheritance further. For now, just consider it a rule that, in all analyzer modules, you should provide this identifier as part of the declaration of analyze.

For those who are knowledgeable about C++, the base class art::EDAnalyzer declares the member function analyze to be pure virtual; so it must be provided by the derived class. The optional member functions of the base class are declared virtual but not pure virutal; do-nothing versions of these member functions are provided by the base class.



In a future version of this documentation suite, more information will be available in the Users Guide in Chapter.

#### 10.7.3.4 The Constructor for the Class First

The next code in the source file (Listing 10.2) is the *definition* of the constructor for the class First. This constructor simply prints some information (via std::cout) to let the user know that it has been called.

```
tex::First::First(fhicl::ParameterSet\ const\&\ pset): art::EDAnalyzer(pset)\ \{std::cout << "Hello\ from\ First::constructor." << std::endl;\ \}
```

The fragment tex::First::First says that this definition is for a constructor of the class First from the namespace tex.

The argument to the constructor is of type fhicl::ParameterSet const& as required by *art*. The class ParameterSet, found in the namespace fhicl, is a C++ representation of a FHiCL parameter set (aka FHiCL *table*). You will learn how to use this parameter set object in Chapter 13.

The argument to the constructor is passed by const reference, const&. This is a requirement specified by *art*; if you write a constructor that does not have exactly the correct argument type, then the compiler will issue a diagnostic and will stop compilation.

The first line of the constructor contains the fragment ": art::EDAnalyzer(pset)". This is the constructor's initializer list and it tells the compiler to call the constructor of the base class art::EDAnalyzer, passing it the parameter set as an argument. This is required by rule 3 in the list in Section 10.7.3.3.

The requirement that the constructor of an analyzer module pass the parameter set to the constructor of art::EDAnalyzer started in *art* version 1.08.09. If you are using an earlier version of *art*, constructors of analyzer modules must NOT call the constructor of art::EDAnalyzer.

#### 10.7.3.5 Aside: Omitting Argument Names in Function Declarations

In the declaration of the class First, you may have noticed that the declaration of the member function analyze supplied a name for its argument (event) but the declaration of the constructor did not supply a name for its argument.

In the declaration of a function, a name supplied for an argument is ignored by the compiler. So code will compile correctly with or without a name. Remember that a constructor is just a special kind of function so the rule applies to constructors too. It is very common for authors of code to provide an argument name as a form of documentation. You will code written both with and without named arguments in declarations.

The above discussion only applied to the *declarations* of functions, not to their definition (aka implementation).

#### 10.7.3.6 The Member Function analyze and the Representation of an Event

The definition of the member function analyze comes next in the source file and is reproduced here

If the type of the argument is not exactly correct, including the the const&, the compiler will issue a diagnostic and stop compilation. The compiler is able to do this because of one of the features of *inheritance*; the details of how this works is beyond the scope of this discussion.

Note that the override contextual identifier that was present in the declaration of this member function is not present in its definition; this is standard C++ usage.

Section 3.6.1 discussed the HEP idea of an event and the *art* idea of a three-part event identifier. The class art::Event is the representation within *art* of the HEP notion of an event. For the present discussion it is safe to consider the following over-simplified view of an event: it contains an event identifier plus a collection of data products (see Section 3.6.4). The name of the argument event has no meaning either to *art* or to the compiler – it is just an identifier – but your code will be easier to read if you choose a meaningful name.



At any given time in a running *art* program there is only ever one art::Event object; in the rest of this paragraph we will call this object *the event*. It is owned and managed by *art*, but *art* lets analyzer modules see the contents of the event; it does so by passing the event by const reference when it calls the analyze member function of analyzer modules. Because the event is passed by reference (indicated by the &), the member function analyze does not get a copy of the event; instead it is told where to find the event. This makes it efficient to pass an event object even if the event contains a lot of information. Because the argument is a const reference, if your code tries to change the contents of the event, the compiler will issue a diagnostic and stop compilation.



As described in Section 3.6.3, analyzer modules may only inspect data in event, not modify it. This section has shown how *art* institutes this policy as a hard rule that will be enforced rigorously by the compiler:

- 1. The compiler will issue an error if an analyzer module does not contain a member function named analyze with exactly the correct signature.
- 2. In the correct signature, the argument event is a const reference.
- 3. Because event is const, the compiler will issue an error if the module tries to call any member function of art::Event that will modify the event.

You can find the header file for art::Event by following the guidelines described in Section 7.6.2. A future version of this documentation will contain a chapter in the Users Guide that provides a complete explanation of art::Event. Here, and in the rest of the Workbook, the features of art::Event will be explained as needed.

The body of the function is almost trivial: it prints some information to let the user know that it has been called. In Section 10.7.1, when you ran *art* using first.fcl, the printout from the first event was

```
Hello from First::analyze. Event id: run: 1 subRun: 0 event: 1
```

If you compare this to the source code you can see that the fragment

```
<< event.id()
```

creates the following printout

```
run: 1 subRun: 0 event: 1
```

This fragment tells the compiler to do the following:

- 1. In the class art::Event, find the member function named id() and call this member function on the object event.
- 2. Whatever is returned by this function call, find its stream insertion operator and call it.

From this description you can probably guess that the member function art::Event::id() returns an object that represents the three-part event identifier. In Section 10.7.3.7 you will learn that this guess is correct.

#### **10.7.3.7 Representing an Event Identifier with art::**EventID

Section 3.6.1 discussed the idea of an event identifier, which has three components, a run number, a subRun number and event number. In this section you will learn where to find the class that *art* uses to represent an event identifier. Rather than simply telling you the answer, this section will guide you through the process of discovering the answer for yourself.

Before you work through this section, you may wish to review Section 7.6 which discusses how to find header files.

In Section 10.7.3.6 you looked at some code and the printout that it made; this strongly suggested that the member function art::Event::id() returns an object that represents the event identifier. To follow up on this suggestion, look at the header file for art::Event. Enter:

#### less \$ART\_INC/art/Framework/Principal/Event.h

or use one of the code browsers discussed in 7.6.2. In this file you will find the definition of the member function id():

```
EventID
id() const {return aux_.id();}
```

The important thing to look at here is the return type, <code>EventID</code>, which looks like a good candidate to be the class that holds the event identifier; you do not need to (or want to) know anything about the data member <code>aux\_</code>. If you look near the beginning of <code>Event.h</code> you will see that it has the line:

```
#include "art/Persistency/Provenance/EventID.h"
```

which looks like a good candidate to be the header file for EventID. Look at this file, e.g.,

#### less \$ART\_INC/art/Persistency/Provenance/EventID.h

and you will discover that it is indeed the header file for EventID; you will also see that

In C++, newlines are treated the same as any other white space; so this could have been written on a single line but the authors of Event.h have adopted a style in which return types are always written on their own line

the class EventID is within the namespace art, making its full name art:: EventID. Near the top of the file you will also see the comments:

```
// An EventID labels an unique readout of the data
// acquisition system, which we call an ``event''.
```

This is another clue that art::EventID is the class we are looking for. Look again at EventID.h; you will see that it has accessor methods that permit you see the three components of the an event identifier:

```
RunNumber_t run() const;
SubRunNumber_t subRun() const;
EventNumber_t event() const;
```

Earlier in EventID.h the C++ type\*\* EventNumber\_t was defined as:

```
namespace art {
  typedef std::uint32_t EventNumber_t;
}
```

meaning that the event number is represented as a 32-bit unsigned integer. A  $typedef(\gamma)$  is a different name, or an alias, by which a type can be identified. If you are not familiar with the C++ concept of typedef, or if you are not familiar with the definite-length integral types defined by the <cstdint> header, consult any standard C++ documentation. If you dig deeper into the layers included in the art::EventID header, you will see that the run number and subRun number are also implemented as 32-bit unsigned integers.



At this point you can be sure that art::EventID is the class that *art* uses to represent the three-part event identifier: the class has the right functionality. It is also true that the comments agree with this hypothesis, but comments are often ill-maintained; be wary of comments and always read the code.

The authors of art might have chosen an alternate definition of EventNumber\_t

```
namespace art {
  typedef unsigned EventNumber_t;
}
```

<sup>\*\*</sup>In C++ the collective noun *type*, refers to both the built-in types, such as int and float, plus user defined types, which include classes, structs and typedefs.

The difference is the use of unsigned rather than std::uint32\_t. This alternate version was not chosen because it runs the risk that some computers might consider this type to have a length of 32 bits while other computers might consider it to have a length of 16 or 64 bits. In the defintion that is used by *art*, an event number is guaranteed to be exactly 32 bits on all computers.

Why did the authors of *art* insert the extra level of indirection and not simply define the following member function inside art::EventID?

```
std::unit32_t event() const;
```

The answer is that it makes it easy to change the definition of the type should that be necessary. If, for example, an experiment requires that event numbers be of length 64 bits, only one change is needed, followed by a recompilation.

It is good practice to use typedefs for every concept for which the underlying data type is not absolutely certain.



It is a very common, but not universal, practice within the HEP C++ community that typedefs that are used to give context-specific names to the C++ built-in types (int, float, char, etc.) end in \_t.

#### 10.7.3.8 DEFINE\_ART\_MACRO: The Module Maker Macros

The final line in First\_module.cc,

```
DEFINE ART MODULE(tex::First)
```

invokes a C preprocessor macro. This macro is defined in the header file that was pulled in by

```
#include "art/Framework/Core/ModuleMacros.h"
```

If you are not familiar with the C preprocessor, don't worry; you do not need to look under the hood. But if you would like to learn about it, consult any standard C++ reference.

The DEFINE\_ART\_MODULE macro instructs the compiler to put some additional code into the shared library made by buildtool. This additional code provides the glue that allows *art* to create instances of the class First without ever seeing the header or the source for the class; it only gets to see the .so and nothing else.



The DEFINE\_ART\_MODULE macro adds two pieces of code to the .so file. It adds a factory function that, when called, will create an instance of First and return a pointer to the base classes art::EDAnalyzer. In this way, art never sees the derived type of any analyzer module; it sees all analyzer modules via pointer to base. When art calls the factory function, it passes as an argument the parameter set specified in the FHiCL file for this module instance. The factory function passes this parameter set through to the constructor of First. The second piece of code put into the .so file is a static object that will be instantiated at load time; when this object is constructed, it will contact the art module registry and register the factory function under the name First. When the FHiCL file says to create a module of type First, art will simply call the registered factory function, passing it the parameter set defined in the FHiCL file. This is the last step in making the connection between the source code of a module and the art instantiation of a module.

#### 10.7.3.9 Some Alternate Styles

C++ allows some flexibility in syntax, which can be seen as either powerful or confusing, depending on your level of expertise. Here we introduce you to a few alternate styles that you will need to recognize and may want to use.

Look at the std::cout line in the analyze method of Listing 10.2:

This could have been written:

This alternate version explicitly creates a temporary object of type art::EventID, whereas the original version created an *implicit* temporary object. When you are first learning C++ it is often useful to break down compound ideas by introducing *explicit tem*-

poraries. However, the recommended best practice is to not introduce explicit temporaries unless there is a good reason to do so.

You will certainly encounter the first line of the above written in a different style, too, i.e.,

```
art::EventID id(event.id());
```

Here id is initialized using *constructor syntax* rather than using *assignment syntax*. For almost all classes these two syntaxes will produce exactly the same result.

You may also see the argument list of the analyze function written a little differently,

```
void analyze( const art::Event& );
instead of
void analyze( art::Event const& );
```

The position of the const has changed. These mean exactly the same thing and the compiler will permit you to use them interchangeably. In most cases, small differences in the placement of the const identifier have very different meanings but, in a few cases, both variants mean the same thing. When C++ allows two different syntaxes that mean the same thing, this documentation suite will point it out.



Finally, Listing 10.3 shows the same information as Listing 10.2 but using a style in which the namespace remains open after the class declaration. In this style, the leading tex: is no longer needed in the definitions of the constructor and of analyze. Both layouts of the code have the same meaning to the compiler. Many experiments use this style in their source code.

Listing 10.3: An alternate layout for First\_module.cc

```
#include "art/Framework/Core/EDAnalyzer.h"
#include "art/Framework/Core/ModuleMacros.h"
#include "art/Framework/Principal/Event.h"

#include <iostream>
namespace tex {

class First : public art::EDAnalyzer {
```

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```
12
     public:
13
14
       explicit First(fhicl::ParameterSet const& );
15
16
       void analyze(art::Event const& event) override;
17
18
     } ;
19
20
     First::First(fhicl::ParameterSet const& pset) : art::EDAnalyzer(pset) {
21
       std::cout << "Hello from First::constructor." << std::endl;</pre>
22
23
24
     void First::analyze(art::Event const& event) {
25
       std::cout << "Hello_from_First::analyze._Event_id:_"</pre>
26
                  << event.id()
27
                  << std::endl;
28
     }
29
30 }
31
32 DEFINE_ART_MODULE(tex::First)
```

## 10.8 What does the Build System Do?

#### 10.8.1 The Basic Operation

In Section 10.4.3 you issued the command buildtool, which *built* First\_module.so. The purpose of this section is to provide some more details about building modules.

When you ran buildtool it performed the following steps:

- 1. It *compiled* First\_module.cc to create an object file (ending in .o).
- 2. It *linked* the object file against the libraries on which it depends and inserted the result into a shared library (ending in . so).

The object file contains the machine code for the class tex::First and the machine code for the additional items created by the DEFINE\_ART\_MODULE C preprocessor macro. The shared library contains the information from the object file plus some additional information that is beyond the scope of this discussion. This process is called *building* the module.

The verb *building* can mean different things, depending on context. Sometimes is just means compiling; sometimes is just means linking; more often, as in this case, it means both.

To be complete, when you ran buildtool it built all of code in the Workbook, both modules and non-modules, but this section will only discuss how it built First\_module.so starting from First\_module.cc.

How did buildtool know what to do? The answer is that it looked in your source directory, where it found a file named CMakeLists.txt; this file contains instructions for cetbuildtools. Yes, when you ran buildtool in your build directory, it did look in your source directory; it knew to do this because, when you sourced setup\_for\_development, it saved the name of the source directory. The instructions in CMakeLists.txt tell cetbuildtools to look for more instructions in the subdirectory ups and in the file art-workbook/CMakeLists.txt, which, in turn, tells it to look for more instructions in the CMakeLists.txt files in each subdirectory of art-workbook.

When **cetbuildtools** has digested these instructions it knows the rules to build everything that it needs to build.

The object file created by the compilation step is a temporary file and, once it has been inserted into the shared library, it is not used any more. Therefore the name of the object file is not important.

On the other hand, the name of the shared library file is very important. art requires that for every module source file (ending in \_module.cc) the build system must create exactly one shared library file (ending in \_module.so). It also requires that the name of each \_module.so file conform to a pattern. Consider the example of the file First\_module.cc; art requires that the shared library for this file match the pattern

```
lib*First_module.so
```

where the \* wildcard matches 0 or more characters.

When naming shared libraries, buildtool uses the following algorithm, which satisfies the *art* requirements and adds some additional features; the algorithm is illustrated using the example of First\_module.cc:

1. find the relative path to the source file, starting from the source directory art-workbook/FirstModule/First module.cc

2. replace all slashes with underscores

```
art-workbook_FirstModule_First_module.cc
```

3. change the trailing .cc to .so
art-workbook\_FirstModule\_First\_module.so

4. add the prefix lib

```
libart-workbook FirstModule First module.so
```

5. put the file into the directory lib, relative to the build directory lib/libart-workbook\_FirstModule\_First\_module.so

You can check that this file is there by issuing the following command from your build directory:

```
Is -I lib/libart-workbook_FirstModule_First_module.so
```

This algorithm guarantees that every module within art-workbook will have a unique name for its shared library.

The experiments using *art* have a variety of build systems. Some of these follow the minimal *art*-conforming pattern, in which the wildcard is replaced with zero characters. If the Workbook had used such a build system, the name of the shared library file would have been

```
lib/libFirst_module.so
```

Both names are legal.

#### 10.8.2 Incremental Builds and Complete Rebuilds

When you edit a file in your source area you will need to rebuild that file in order for those changes to take effect. If any other files in your source area depend on the file that you edited, they too will need to be rebuilt. To do this, reissue the command:

#### buildtool

Remember that this command must be executed from your build directory and that, before executing it, you must have setup the environment in your build window. When you run this command, **cetbuildtools** will automatically determine which files need to be rebuilt and will rebuild them; it will not waste time rebuilding files that do not need to be rebuilt. This is called an *incremental build* and it will usually complete much faster than the initial build.

If you want to clean up everything in your build area and rebuild everything from scratch, use the following command:

#### buildtool -c

This command will give you five seconds to abort it before it starts removing files; to abort, type ctrl-C in your build window. It will take about the same time to execute as did your initial build of the Workbook. The name of the option -c is a mnemonic for "clean".

When you do a clean build it will remove all files in your build directory that are not managed by **cetbuildtools**. For example, if you redirected the output of *art* as follows,

#### art -c fcl/FirstModule/first.fcl >& first.log



then, when you do a clean build, the file first.log will be deleted. This is why the instructions earlier in this chapter told you to redirect outtut to a log file by

### art -c fcl/FirstModule/first.fcl >& output/first.log

When you ran buildtool, it created a directory to hold your output files and you created a symbolic link, named output, from your build directory to this new directory. Both the other directory and the symbolic link survive clean builds and your output files will be preserved. The Workbook exercises write all of their root and event-data output files to this directory.

If you edit certain files in the ups subdirectory of your source directory, rebuilding requires an extra step. If you edit one of these files, the next time that you run buildtool, it will issue an error message saying that you need to re-source setup\_for\_development. If you get this message, make sure that you are in your build directory, and

 $source \ ../art-workbook/ups/setup\_for\_development \ \backslash -p \ \$ART\_WORKBOOK\_QUAL$ 

buildtool

#### 10.8.3 Finding Header Files at Compile Time

When setup\_for\_development establishes the working environment for the build directory, it does a UPS setup on the UPS products that it requires; this triggers a chain of additional UPS setups. As each UPS product is set up, that product defines many environment variables, two of which are <PRODUCT-NAME>\_INC and <PRODUCT-NAME>\_LIB. The first of these points to a directory that is the root of the header file hierarchy for that version of that UPS product. The second of these points to a single directory that holds all of the shared library files for that UPS product.

You can spot-check this by doing, for example,

```
Is $TOYEXPERIMENT_INC/*
Is $TOYEXPERIMENT_LIB
Is $ART_INC/*
Is $ART_LIB
```

You will see that the \_INC directories have a subdirectory tree underneath them while the \_LIB directories do not.

There are a few small perturbations on this pattern. The most visible is that the ROOT product puts most of its header files into a single directory, \$ROOT\_INC and does not clone the directory hierarchy of its source files. The Geant4 product does the same thing.

When the compiler compiles a .cc file, it needs to know where to find the files specified by the #include directives. The compiler looks for included files by first looking for arguments on the command line of the form

```
-I<absolute-path-to-a-directory>
```

There may be many such arguments on one command line. The compiler assembles the set of all -l arguments and uses it as an include path; that is, it looks for the header files by trying the first directory in the path and if it does not find it there, it tries the second directory in the path, and so on. The choice of -l for the name of the argument is a mnemonic for Include.

When buildtool compiles a .cc file it adds many -l options to the command

line; it adds one for each UPS product that was set up when you sourced setup\_for\_development. When building First\_module.cc, buildtool added - I\$ART\_INC, -I\$TOYEXPERIMENT\_INC and many more.

A corollary of this discussion is that when you wish to include a header file from a UPS product, the #include directive must contain the relative path to the desired file, starting from the \_INC environment variable for that UPS product.



This system illustrates how the Workbook can work the same way on many different computers at many different sites. As the author of some code, you only need to know paths of include files relative to the relevant \_INC environment variable. This environment variable may have different values from one computer to another but the setup and build systems will ensure that the site-specific information is communicated to the compiler using environment variables and the -I option.

This system has the potential weakness that if two products each have a header file with exactly the same relative path name, the compiler will get confused. Should this happen, the compiler will always choose the file from the earlier of the two-l arguments on the command line, even when the author of the code intended the second choice to be used. To mitgate this problem, the *art* and UPS teams have adopted the convention that, whenever possible, the first element of the relative path in an #include directive will be the UPS package name. It is the implementation of this convention that led to the repeated directory name art-workbook/art-workbook that you saw in your source directory. There are a handful of UPS products for which this pattern is not followed and they will be pointed out as they are encountered.

The convention of having the UPS product name in the relative path of #include directives also tells readers of the code where to look for the included file.

## 10.8.4 Finding Shared Library Files at Link Time

The module First\_module.cc needs to call methods of the class art::Event. Therefore the compiler left a notation in the object file saying "to use this object file you need to tell it where to find art::Event." The technical way to say this is that the object file contains a list of *undefined symbols* or *undefined external references*. When the linker makes the shared library

libart-workbook\_FirstModule\_First\_module.so

it must resolve all of the undefined symbols from all of the object files that go into the library. To resolve a symbol, the linker must learn what shared library defines that symbol. When it discovers the answer, it will write the name of that shared library into something called the *dependency list* that is kept inside the shared library. **cetbuildtools** tells the linker that the dependency list should contain only the filename of each shared library, not the full path to it. If, after the linker has finished, there remain unresolved symbols, then the linker will issue an error message and the build will fail.

Dependency lists are not recursive. If library A depends on library B and library B depends on library C, then the dependency list of library A should contain only library B. In other words, the dependency list is allowed to contain only *direct dependencies* (also called *first order* dependencies).

To learn where to look for symbol definitions, the linker looks at its command line to find something called the *link list*. The link list can be specified in several different ways and the way that **cetbuildtools** uses is simply to write the link list as the absolute path to every .so file that the linker needs to know about. The link list can be different for every shared library that the build system builds. However it is very frequently true that if a directory contains several modules, then all of the modules will require the same link list. The bottom line is that the author of a module needs to know the link list that is needed to build the shared library for that module.



For these Workbook exercises, the author of each exercise has determined the link list for each shared library that will be built for that exercise. In the <code>cetbuildtools</code> system, the link list for <code>First\_module.cc</code> is located in the <code>CMakeLists.txt</code> file from same directory as <code>First\_module.cc</code>; the contents of this file are shown in Listing 10.4. This <code>CMakeLists.txt</code> file says that all modules found in this directory should be built with the same link list and it gives the link list; the link list is the seven lines that begin with a dollar sign; these lines each contain one cmake variable. Recall that <code>cetbuildtools</code> is a build system that lives on top of cmake, which is another build system. A cmake variable is much like an environment variable except that is only defined within the environment of the running build system; you cannot look at it with printeny.

The five cmake variables beginning with ART\_ were defined when buildtool set up the UPS *art* product. Each of these variables defines an absolute path to a shared library in

Part

```
$ART_LIB. For example ${ART_FRAMEWORK_CORE} resolves to
```

```
$ART_LIB/libart_Framework_Core.so
```

Almost all *art* modules will depend on these five libraries. Similarly the other two variables resolve to shared libraries in the **fhiclcpp** and **cetlib** UPS products.

When **cetbuildtools** constructs the command line to run the linker, it copies the link list from the CMakeLists.txt file to the command linker line.

The experiments that use *art* use a variety of build systems. Some of these build systems do not require that all external symbols be resolved at link time; they allow some external symbols to be resolved at run-time. This is legal but it can lead to certain difficulties. A future version of this documentation suite will contain a chapter in the Users Guide that discusses linkage loops and how use of closed links can prevent them. This section will then just reference it.

Consult the cmake and **cetbuildtools** documentation to understand the remaining details of this file.

Listing 10.4: The file art-workbook/FirstModule/CMakeLists.txt

```
1
  art_make (MODULE_LIBRARIES
2
    ${ART_FRAMEWORK_CORE}
3
    ${ART_FRAMEWORK_PRINCIPAL}
4
    ${ART_PERSISTENCY_COMMON}
5
    ${ART_FRAMEWORK_SERVICES_REGISTRY}
6
    ${ART_FRAMEWORK_SERVICES_OPTIONAL}
7
    ${FHICLCPP}
8
    ${CETLIB}
```

## 10.8.5 Build System Details

This section provides the next layer of details about the build system; in a future version of this documentation set, the Users Guide will have a chapter with all of the details. This entire section contains expert material.



If you want to see what buildtool is actually doing, you can enable verbose mode by issuing the command:

buildtool VERBOSE=TRUE

Step	Flags
Compiler	-Dart_workbook_FirstModule_First_module_EXPORTS
	-DNDEBUG
Linker	-Wl,-no-undefined -shared
Both	-O3 -g -fno-omit-frame-pointer -Werror -pedantic
	-Wall -Werror=return-type -Wextra -Wno-long-long -Winit-self
	-Woverloaded-virtual -std=c++11
	-D_GLIBCXX_USE_NANOSLEEP -fPIC

Table 10.1: Compiler and Linker Flags for a Profile Build

For example, if you really want to know the name of the object file, you can find it in the verbose output. For this exercise, the object file is

```
./art-workbook/FirstModule/CMakeFiles/
art-workbook_FirstModule_First_module.dir/First_module.cc.o
```

where the above is really just one line.

Also, you can read the verbose listing to discover the flags given to the compiler and linker. The more instructive compiler and linker flags valid at time of writing are given in Table 10.1. The C++ 11 features are selected by the presence of the -std=c++11 flag and a high level of error checking is specified. The linker flag,

```
-Wl, --no-undefined
```

tells the linker that it must resolve all external references at link time. This is sometime referred to as a *closed link*.

## 10.9 Suggested Activities

This section contains some suggested exercises in which you will make your own modules and learn more about how to use the class art::EventID.

## 10.9.1 Create Your Second Module

In this exercise you will create a new module by copying First\_module.cc and making the necessary changes; you will build it using buildtool; you will copy first.fcl

and make the necessary changes; and you will run the new module using the new FHiCL file.

Go to your source window and cd to your source directory. If you have logged out, out remember to re-establish your working environment; see Section 10.5 Type the following commands:

cd art-workbook/FirstModule

cp First\_module.cc Second\_module.cc

cp first.fcl second.fcl

Edit the files Second\_module.cc and second.fcl. In both files, change every occurence of the string "First" to "Second"; there are eight places in the source file and two in the FHiCL file, one of which is in a comment.

The new module needs the same link list as did First\_module.cc so there is no need to edit CMakeLists.txt; the instructions in CMakeLists.txt tell buildtool to build all modules that it finds in this directory and to use the same link list for all modules.

Go to your build window and cd to your build directory. Again, remember to re-establish your working environment as necessary. Rebuild the Workbook code:

### buildtool

This should complete with the message:

INFO: Stage build successful.

If you get an error message, consult a local expert or the *art* team as described in Section 3.4.

When you run buildtool it will perform an incremental build (see Section 10.8.2) during which it will detect Second\_module.cc and build it.

You can verify that buildtool created the expected shared library:

Is lib/\*Second\*.so

```
lib/libart-workbook_FirstModule_Second_module.so
```

Stay in your build directory and run the new module:

```
art -c fcl/FirstModule/second.fcl >& output/second.log
```

Compare output/second.log with output/first.log. You should see that "First" has been replaced by "Second" everywhere and the date/time lines are different.

#### 10.9.2 Use artmod to Create Your Third Module

This exercise is much like the previous one; the difference is that you will use a tool named artmod to create the source file for the module.

Go to your source window and cd to your source directory. If you have logged out, remember to re-establish your working environment; see Section 10.5

The command artmod creates a file containing the skeleton of a module. It is supplied by the UPS product **cetpkgsupport**, which was set up when you performed the last step of establishing the environment in the source window, sourcing setup\_deps. You can verify that the command is in your path by using the bash built-in command type (output shown on two lines):

#### type artmod

```
artmod is hashed (/ds50/app/products/cetpkgsupport/ v1_02_00/bin/artmod)
```

The leading elements of the directory name will reflect your UPS products area, and may be different from what is shown here.

From your source directory, type the following commands:

```
cd art-workbook/FirstModule
artmod analyzer tex::Third
cp first.fcl third.fcl
```

The second command tells art mod to create a source file named Third\_module.cc that contains the skeleton for an 'analyzer' module, to be named Third in the namespace

tex.

If you compare Third\_module.cc to First\_module.cc you will see a few differences:

- 1. Third\_module.cc is longer and has more comments
- 2. the layout of the class is a little different but the two layouts are equivalent
- 3. there are some extra #include directives
- 4. the include for <iostream> is missing
- 5. a argument (p) is supplied in the definition of the constructor
- 6. in the analyze member function, the name of the argument is different (event vs e)
- 7. artmod supplies the skeleton of a destructor ( $\sim$ Third)

The #include directives provided by artmod are a best guess, made by the author of artmod, about which ones will be needed in a "typical" module. Other than slowing down the compiler by an amount you won't notice, the extra #include directives do no harm; keep them or leave them as you see fit.

Edit Third module.cc

- 1. add the #include directive for <iostream>
- 2. copy the bodies of the constructor and the analyze member function from First\_module.cc; change the string "First" to "Third"
- 3. delete the argument from the definition of the constructor
- 4. in the definition of the member function analyze, change the name of the argument to event.

When you built First\_module.cc, the compiler wrote a destructor for you that is identical to the destructor written by artmod; so you can leave the destructor as artmod wrote it, i.e., with an empty body. This class has no work to do in the destructor.

Edit third.fcl Change every occurrence of the string "First" to "Third"; there are two places, one of which is in a comment.

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Go to your build window and cd to your build directory. If you have logged, out remember to re-establish your working environment; see Section 10.5. Rebuild the Workbook code:

#### buildtool

Refer to the previous section to learn how to identify a successful build and how to verify that the expected library was created.

Stay in your build directory and run the third module:

```
art -c fcl/FirstModule/third.fcl >& output/third.log
```

Compare output/third.log with output/first.log. You should see that the printout from First\_module.cc has been replaced by that from Third\_module.cc. artmod has many options that you can explore by typing:

```
artmod -help
```

## 10.9.3 Running Many Modules at Once

In this exercise you will run four modules at once, the three made in this exercise plus the HelloWorld module from Chapter 9.

Go to your source window and cd to your source directory. Type the following commands:

```
cd art-workbook/FirstModule
```

```
cp first.fcl all.fcl
```

Edit the file all.fcl and replace the physics parameter set with the contents of Listing 10.5. This parameter set:

- 1. defines four module labels and
- 2. puts all four module labels into the end\_paths sequence.

When you run *art* on this FHiCL file, *art* will first look at the definition of end\_paths and learn that you want it to run four module labels. Then it will look in the analyzers parameter set to find the definition of each module label; in each definition *art* will find

the class name of the module that it should run. Given the class name and the environment variable LD\_LIBRARY\_PATH, *art* can find the right shared library to load. If you need a refresher on module labels and end\_paths, refer to Sections 9.8.7 and 9.8.8.

Listing 10.5: The physics parameter set for all.fcl

```
1
  physics :{
2
    analyzers: {
      hello : {
3
4
         module_type : HelloWorld
5
6
       first : {
7
         module_type : First
8
9
       second : {
10
        module_type : Second
11
12
       third : {
13
         module_type : Third
14
15
    }
16
17
          : [ hello, first, second, third ]
18
     end_paths : [ e1 ]
19
20
```

Go to your build window and cd to your build directory. If you have logged, out remember to re-establish your working environment; see Section 10.5. You do not need to build any code for this exercise.

Run the exercise:

## art -c fcl/FirstModule/all.fcl >& output/all.log

Compare output/all.log with the log files from the previous exercises. The new log file should contain printout from each of the four modules. Once, near the start of the file, you should see the printout from the three constructors; remember that the HelloWorld module does not make any printout in its constructor. For each event you should see the printout from the four analyze member functions.

Remember that *art* is free to run analyzer modules in any order; this was discussed in Section 9.8.8.

#### 10.9.4 Access Parts of the EventID

In this exercise, you will access the individual parts of the event identifier.

Before proceeding with this section, review the material in Section 10.7.3.7 which discusses the class art::EventID. The header file for this class is:

```
$ART_INC/art/Persistency/Provenance/EventID.h
```

In this exercise, you are asked to rewrite the file Second\_module.cc so that the printout made by the analyze method looks like the following (lines split here due to space restrictions):

```
Hello from FirstAnswer01::analyze. run number: 1
    sub run number: 0 event number: 1
Hello from FirstAnswer01::analyze. run number: 1
    sub run number: 0 event number: 2
```

and so on for each event.

To do this, you will need to reformat the text in the std::cout statement and you will need to separately extract the run, subRun and event numbers from the art::EventID object.

You will do the editing in your source window, in the subdirectory art-workbook/ FirstModule.

When you think that you have successfully rewritten the module, you can test it by going to your build window and cd'ing to your build directory. Then:

#### buildtool

```
art -c fcl/FirstModule/second.fcl >& output/eventid.log
```

If you have not figured out how to do this exercise after about 15 minutes, you can find one possible answer in the file FirstAnswer01\_module.cc, in the same directory as First module.cc.

To run the answer module and verify that it makes the requested output, run (command can be typed on a single line):

```
art -c fcl/FirstModule/firstAnswer01.fcl >& \output/firstAnswer01.log
```

You did not need to build this module because it was already built the first time that you ran buildtool; that run of buildtool built all of the modules in the Workbook.

There is a second correct answer to this exercise. If you look at the header file for art:: Event, you will see that this class also has member functions

```
EventNumber_t event() const {return aux_.event();}
SubRunNumber_t subRun() const {return aux_.subRun();}
RunNumber_t run() const {return id().run();}
```

So you could have called these directly,

instead of

But the point of this exercise was to learn a little about how to dig down into nested header files to find the information you need.

## 10.10 Final Remarks

## 10.10.1 Why is there no First\_module.h File?

When you performed the exercises in this chapter, you saw, for example, the file First\_module.cc but there was no corresponding First\_module.h file. This section will explain why.

In a typical C++ programming environment there is a header file (.h) for each source file (.cc). As an example, consider the files Point.h and Point.cc that you saw in

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#### Section 6.6.10.

The reason for having Point.h is that the implementation of the class, Point.cc, and the users of the class need to agree on what the class Point is. In the Section 6.6.10 example, the only user of the class is the main program, ptest.cc. The file Point.h serves as the unique, authoritative declaration of what the class is; both Point.cc and ptest.cc rely on on this declaration.

If you think carefully, you are already aware of a very common exception to the pattern of one .h file for each .cc file: there is never a header file for a main program. For example, in the examples that exercised the class Point, ptest.cc had no header file. Why not? No other piece of user-written code needs to know about any classes or functions declared or defined inside ptest.cc.

The First\_module.h file is omitted simply because every entity that needs to see the declaration of the class First is already inside the file First\_module.cc. There is no reason to have a separate header file. Recall the "dangerous bend" paragraph at the end of Section 10.7.3.8 that described how *art* is able to use modules without needing to know about the declaration of the module class.

art is designed such that only art may construct instances of module classes and only art may call member functions of module classes. In particular, modules may not construct other modules and may not call member functions of other modules. The absence of a First\_module.h, provides a physical barrier that enforces this design.

#### 10.10.2 The Three-File Module Style

In this chapter, the source for the module First was written in a single file. You may also write it using three files, First.h, First.cc and First\_module.cc.



Some experiments use this three-file style. The authors of *art* do not recommend it, however, because it exposes the declaration of First in a way that permits it to be misused (as was discussed in Section 10.10.1). The build system distributed with the Workbook has not been configured to build modules written in this style.

In this style, First.h contains the class declaration plus any necessary #include directives; it now also requires code guards (see Section 32.8); this is shown in Listing 10.6.

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Listing 10.6: The contents of First.h in the three-file model

```
1
   #ifndef art-workbook_FirstModule_First_h
   #define art-workbook_FirstModule_First_h
3
4
5
  #include "art/Framework/Core/EDAnalyzer.h"
   #include "art/Framework/Principal/Event.h"
7
8
  namespace tex {
9
10
     class First : public art::EDAnalyzer {
11
12
     public:
13
14
       explicit First(fhicl::ParameterSet const& );
15
16
       void analyze(art::Event const& event) override;
17
18
     } ;
19
20
21
  #endif
```

The file First.cc contains the definitions of the constructor and the analyze member function, plus the necessary #include directives; this is shown in Listing 10.7.

Listing 10.7: The contents of First.cc in the three-file model

```
1
   #include "art-workbook/FirstModule/First.h"
3
4
   #include <iostream>
5
6
  tex::First::First(fhicl::ParameterSet const& pset ) : art::EDAnalyzer(pset) {
7
     std::cout << "Hello_from_First::constructor." << std::endl;</pre>
8
   }
9
10
   void tex::First::analyze(art::Event const& event) {
     std::cout << "Hello_from_First::analyze._Event_id:_"</pre>
11
12
                << event.id()
13
                << std::endl;
14
```

And First\_module.cc is now stripped down to the invocation of the DEFINE\_ART\_MODULE macro plus the necessary #include directives; this is shown in Listing 10.8.

Listing 10.8: The contents of First\_module.cc in the three-file model

```
1
2 #include "art-workbook/FirstModule/First.h"
3 #include "art/Framework/Core/ModuleMacros.h"
4
5 DEFINE_ART_MODULE(tex::First)
```

## 10.11 Flow of Execution from Source to FHiCL File

The properties that a class must have in order to be an analyer module are summarized in Section 10.7.3.2 for reference. This section reviews how the source code found in an analyzer module, e.g., First\_module.cc, is executed by *art*:

- 1. The script setup\_for\_development defines many environment variables that are used by buildtool, *art* and toyExperiment.
- 2. LD\_LIBRARY\_PATH, an important environment variable, contains the directory lib in your build area plus the lib directories from many UPS products, including *art*.
- 3. buildtool compiles First\_module.cc to a temporary object file.
- 4. buildtool links the temporary object file to create a shared library in the lib subdirectory of your build area:

```
lib/libart-workbook_FirstModule_First_module.so
```

- 5. When you run *art* using file first.fcl, this file tells *art* to find and load a module with the "module\_type" First.
- 6. In response to this request, *art* will search the directories in LD\_LIBRARY\_PATH to find a shared library file whose name matches the pattern:

```
lib*First_module.so
```

- 7. If *art* finds either zero or more than one match to this pattern, it will issue an error message and stop.
- 8. If art finds exactly one match to this pattern, it will load the shared library.
- 9. After *art* has loaded the shared library, it has access to a function that can, on demand, create instances of the class First.

The last bullet really means that the shared library contains a factory function that can construct instances of First and return a pointer to the base class, art::EDANalyzer. The shared library also contains a static object that, at load-time, will contact the *art* module registry and register the factory function under the module\_type First.



## 11 Keeping Up to Date with Workbook Code and Documentation

#### 11,1 Introduction

As you well know by now, the Workbook exercises require you to download some code to edit, build, execute and evaluate. Both the documentation and the code it references are expected to undergo continual development throughout 2014. The latest is always available at the art Documentation website.



Announcements of new releases are made on the art-users@fnal.gov mailing list.

If you are continuing directly from Exercise 1 to Exercise 2, you have the choice of updating to a newer release if one exists, or you can just continue on with the release you have until you finish the exercises available in it. At that point, you will need to update to a newer release (with more exercises) to continue. Come back to this chapter whenever you reach the end of the available exercises. Or come back and update whenever a new release is announced; it may include improvements to existing exercises.

#### How to Update 11.2

In order to update, you need to:

- 1. Determine whether an updated release is available, and what release it is.
- 2. Switch to the updated documentation.

Part

- 3. In your source window, use git to update your working version of the code in the (higher-level) art-workbook directory
- 4. In your build window, build the new version of the code.

## 11.2.1 Get Updated Documentation

First, check which documentation release you're currently using: it's noted on the title page\*. Then go to the *art* Documentation website and compare your documentation release number to the latest available.

Download a new copy of the documentation, as needed.

#### 11.2.2 Get Updated Code and Build It

Also noted on the title page of the documentation is the release<sup>†</sup> of the art-workbook code that the documentation is intended for. Recall from Figure 10.1 that git commands are used to clone the code in the remote repository into your local copy, then copy the requested release from that local copy into your working area. The git system is described in more detail in Chapter 22.

Chances are that you're using the code release that goes with the documentation you have been using. You can check by looking in the file art-workbook/ups/product\_deps. From your source directory run:

```
grep art_workbook ups/product_deps
```

```
parent art_workbook v0_00_13
```

This shows version v0\_00\_13 as an example. If your version is earlier than the one listed on the cover of the latest documentation, you will need to get new code and build it.

These instructions illustrate updating the working version of the art-workbook code from version v0\_00\_13 to version v0\_00\_15. There is nothing special about these two versions; the instructions serve as a model for a change between any pair of versions.

<sup>\*</sup>Versions of the *art* documentation prior to 0.70 do not have this information on the front page; for these versions, the required version of art-workbook can be found in the section "Setting up to Run Exercises" in Exercise 2.

<sup>&</sup>lt;sup>†</sup>The terms "release" and "version" are used interchangeably here.

- 1. Start from (or cd to) your source directory (see Section 10.4.1).
- 2. Use git status and make a note of the files that you have modified and/or added (see Section 22.1.3 for instructions).
- 3. Switch from your tagged version branch back to the develop branch ("branches" are discussed in Chapter 22, you don't need to understand them at this stage<sup>‡</sup>).

## git checkout develop

```
Switched to branch 'develop'
```

4. Update your local copy of the respository (the .git directory)

```
git pull
```

The output from this command is shown in Listing 11.1.

5. Switch your working code to the new branch:

```
git checkout -b v0_00_15 v0_00_15

Switched to a new branch 'v0_00_15'
```

In the messages produced in this step, watch for the names of files that you have modified. Check for conflicts that git did not merge correctly.

To rebuild your updated working code:

- 1. In your build window, cd to your build directory
- 2. Tell **cetbuildtools** to look for, and act on, any changes in your checked out version of the code (command shown on two lines):

```
source ../art-workbook/ups/setup_for_development \-p $ART_WORKBOOK_QUAL
```

3. Rebuild:

buildtool

If this step does not complete successfully, the first thing to try is a clean rebuild:

<sup>‡</sup>If you are familiar with git concepts, you may want to know this: The authors of the art-workbook follow the convention that they make a new git branch for every release of art-workbook and the name of the branch matches the version number of art-workbook. In the current example, the local working environment knows about two branches, the develop branch and the branch for version v0\_00\_13. The develop branch is the name of the branch that always contains the most recent art-workbook code.

**Listing 11.1:** Example of the output produced by git pull

```
1 From http://cdcvs.fnal.gov/projects/art-workbook
      e79d9ef..81d2a76 develop -> origin/develop 6435ecc..c0c1af5 master -> origin/master
2
3
4 From http://cdcvs.fnal.gov/projects/art-workbook
5 * [new tag] v0_00_14 -> v0_00_14
6 * [new tag] v0_00_15 -> v0_00_15
7
  Updating e79d9ef..81d2a76
8 Fast-forward
   art-workbook/ModuleInstances/magic.fcl | 26 +++++++++++------
  art-workbook/ParameterSets/PSet01_module.cc | 36 ++++++++++++++++-----
10
11
    art-workbook/ParameterSets/PSet02 module.cc | 53 +++++++++++++++------
    art-workbook/ParameterSets/PSet03_module.cc | 28 +++++++++++-----
12
13
    art-workbook/ParameterSets/PSet04_module.cc | 44 ++++++++++++++++-----
    art-workbook/ParameterSets/pset01.fcl
14
                                                | 6 ++---
15
    art-workbook/ParameterSets/pset02.fcl
                                               | 14 +++++----
16
    art-workbook/ParameterSets/pset03.fcl
    art-workbook/ParameterSets/pset04.fcl
                                                7 +++---
17
18
    ups/product_deps
                                                 | 2 +-
19
   10 files changed, 109 insertions (+), 113 deletions (-)
```

buildtool -c

## 11.2.3 See which Files you have Modified or Added

At any time you can check to see which files you have modified and which you have added. The code is structured in such a way that when you checkout a new version, these files will remain in your working directory and will not be modified or deleted. The git checkout command will generate some informational messages about them, but you do not need to take any action.

To see the new/modified files, cd to your source directory and issue the git status command. Suppose that you have checked out version v0\_00\_13, modified first.fcl and added second.fcl. The git status command will produce the following output:

#### git status

```
# On branch v0_00_13
# Changes not staged for commit:
```

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```
# (use "git add <file>..." to update what will be committed)
# (use "git checkout -- <file>..." to discard changes in
working directory)
#
# modified: first.fcl
#
# Untracked files:
# (use "git add <file>..." to include in what will be committed)
#
# second.fcl
no changes added to commit (use "git add" and/or "git commit -a")
```



Do not issue the git add or git commit commands that are suggested in the command output above.

In the rare case that you have neither modified nor added any files, the output of git status will look like:

## git status

```
# On branch v0_00_13
```

# 12 Exercise 3: Some other Member Functions of Modules

## 12.1 Introduction

Recall the discussion in Section 3.6.2 about widget-making workers on an assembly line. All workers have a task to perform on each widget as it passes by and some workers may also need to perform start-up or shut-down tasks. If a module has something that it must do at the start of the job, then the author of the module can write a member function named beginJob() that performs these tasks. Similarly the author of a module can write a member function named endJob to do tasks that need to be performed at the end of the job. *art* will call both of these member functions at the appropriate time.

The author of a module may also provide member functions to perform actions at the start of a subRun, at start of a run, at the end of a subRun or at the end of a Run.

These member functions are *optional*; i.e., they are always allowed in a module but never required. They have prescribed names and argument lists.



In this exercise you will build and execute an analyzer module that illustrates three of these member functions: beginJob, beginRun and beginSubRun. These member functions are called, respectively, once at the start of the *art* job, once for each new run and once for each new subRun.

You may also perform a suggested exercise to add the three corresponding member functions endJob, endRun and endSubRun.

## 12.2 Prerequisites

The prerequisites for this chapter include all of the material in Part I (Introduction) and all of the material up to this point in Part II (Workbook).

In particular, make sure that you understand the *event loop* (see Section 3.6.2).

## 12.3 What You Will Learn

This chapter will show you *how* to provide the optional member functions in your *art* modules to execute special functionality at the beginning and end of jobs, runs and/or subRuns. These include

```
    beginJob()
    beginRun(art::Run const&)
    beginSubRun(art::SubRun const&)
    endJob()
    endRun(art::Run const&)
    endSubRun(art::SubRun const&)
```

As you gain experience, you will gain proficiency at knowing when to provide them.

You will also be introduced to the classes

```
    art::RunID
    art::Run
    art::SubRunID
    art::SubRun
```

that are analogous to the art::EventID and art::Event classes that you have already encountered.

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## 12.4 Setting up to Run this Exercise

Follow the instructions in Section 10.5 if you are logging in after having closed an earlier session. If you are continuing on directly from the previous exercise, keep both your source and build windows open.

## 12.5 Files Used in this Exercise

In your source window, look at the contents of the directory for this exercise, called OptionalMethods:

## Is art-workbook/OptionalMethods

```
CMakeLists.txt OptionalAnswer01_module.cc Optional_module.cc optionalAnswer01.fcl optional.fcl
```

The source code for the module you will run is Optional\_module.cc and the FHiCL file to run it is optional.fcl. The file CMakeLists.txt is identical to that used by the previous exericse since the new features introduced by this module do not require any modifications to the link list. The other two files relate to the exercise you will be asked to do in Section 12.9.

In your build window, just make sure that you are in your build directory. All the code for this exercise is already built; this happened the first time that your ran buildtool.

## 12.6 The Source File Optional\_module.cc

In your source window, look at the source file Optional\_module.cc and compare it to First\_module.cc. The differences are

- 1. it has two new include directives, for Run . h and SubRun . h
- 2. the name of the class has changed from First to Optional
- 3. the Optional class declaration declares three new member functions void beginJob () override;

```
void beginRun ( art::Run const& run ) override;
void beginSubRun( art::SubRun const& subRun ) override;
```

- 4. the text printed by the constructor and analyze member functions has changed
- 5. the file contains the definitions of the three new member functions, each of which simply makes some identifying printout

## 12.6.1 About the begin\* Member Functions

The optional member functions beginJob, beginRun and beginSubRun, described in the Introduction to this chapter (Section 12.1), must have exactly the argument list prescribed by *art* as shown in list item 3 above.

art knows to call the beginJob member function of each module, if present, once at the start of the job; it knows to call beginRun, if present, at the start of each run and, likewise, beginSubRun at the start of each subRun.

#### 12.6.2 About the art::\*ID Classes

In Section 10.7.3.7 you learned about the class art::EventID, which describes the three-part event identifier. *art* also provides two related classes:

```
o art::RunID, a one-part identifier for a run number
```

o art::SubRunID, a two-part identifier for a subRun

The header files for these classes are found at:

```
$ART_INC/art/Persistency/Provenance/RunID.h
$ART_INC/art/Persistency/Provenance/SubRunID.h
```

Similar to the art::Event class discussed in Section 10.7.3.6, art provides art::Run and art::subRun. These contain the IDs, e.g., art::RunID, plus the data products for the entire run or subRun, respectively. You can find their header files at:

```
$ART_INC/art/Framework/Principal/Run.h
$ART_INC/art/Framework/Principal/SubRun.h
```

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In the call to beginSubRun the argument is of type art::SubRun const&. A simplified description of this object is that it contains an art::SubRunID plus a collection of data products that describe the subRun. All of the comments about the class art::Run in the preceding few paragraphs apply to art::SubRun. You can find the header file for art::SubRun at:

less \$ART\_INC/art/Framework/Principal/SubRun.h

#### 12.6.3 Use of the override Identifier

The override identifier on each of these member functions instructs the compiler to check that both the name (and spelling) of the member function and its argument list are correct; if not, the compiler will issue an error message and stop. This is a very handy feature. Without it, a misspelled function name or incorrect argument list would cause the compiler to assume that you intended to define a *new* member function unrelated to one of these optional *art*-defined member functions. This would result in a difficult-to-diagnose run-time error: *art* would simply not recognize your member function and would never call it.

Always provide the override identifier when using any of the optional *art*-defined member functions.



For those with some C++ background, the three member functions beginJob, beginRun and beginSubRun are declared as virtual in the base class, art::EDAnalyzer. The override identifier is new in C++-11 and will not be described in older text books. It instructs the compiler that this member function is intended to override a virtual function from the base class; if the compiler cannot find such a function in the base class, it will issue an error.



#### 12.6.4 Use of const References

In Optional\_module.cc the argument to the beginRun member function is a const reference to an object of type art::Run that holds the current run ID and the collection of data products that together describe the run. If you take a snapshot of a running *art* job you will see that, at any time, there is exactly one object of type art::Run. This object is

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owned by art. art gives modules access to it when it (art) calls the modules' beginRun and endRun member functions.

Because the object is passed by reference, the beginRun member function does not get a copy of the object; instead it is given access to it. Because it is passed by const reference in this example, your analyzer module may look at information in the object but it may not add or change information to the art::Run object.

There is a very important habit that you need to develop as a user of *art*. Many member functions in *art*, in the Workbook code and very likely in your experiment's code, will return information by & or by const&. If you receive these by value, not by reference, then you will make copies that waste both CPU and memory; in some cases these can be significant wastes. Unfortunately there is no way to tell the compiler to catch this mistake. The only solution is your own vigilance.



To access the art::Run and art::SubRun objects through, for example, an art::Event named event, you can use

```
art::SubRun const& subRun = event.getSubRun();
for the subRun and
  art::Run const& run = subRun.getRun();
for the run.
```

## **12.6.5** The analyze Member Function

In your analyze member function, if you have an art::Event, named event, you can access the associated run information by:

```
art::Run const& run = event.getRun();
```

You may sometimes see this written as:

```
auto const& run = event.getRun();
```



Both versions mean exactly the same thing. When a type is long and awkward to write, the auto identifier is very useful; however it is likely to be very confusing to beginners. When you encounter it, check the header files for the classes on the right hand side of the

Part

assignment; from there you can learn the return type of the member function that returned the information.

## 12.7 Running this Exercise

Look at the file optional.fcl. This FHiCL file runs the module Optional on the the input file inputFiles/input03\_data.root. Consult Table 9.1 and you will see that this file contains 15 events, all from run 3. It contains events 1 through 5 from each of subRuns 0, 1 and 2. With this knowledge, and the knowledge of the source file Optional\_module.cc, you should have a clear idea of what this module will print out.

In your build directory, run the following command

## art -c fcl/OptionalMethods/optional.fcl >& output/optional.log

The part of the printed output that comes from the module Optional is given in Listing 12.1. Is this what you expected to see? If not, understand why this module made the printout that it did. If you did not get this printout, double check that you followed the instructions carefully; if that still does not fix it, ask for help (see Section 3.4).

## **12.8** The Member Function beginJob versus the Constructor

The member function beginJob gets called once at the start of the job. The constructor of the each module is also called once at the start of the job. This brings up the question: What code belongs in the constructor and what code belongs in the beginJob member function?

art requires that some tasks be done in the constructor. You have not yet encountered them; they will be pointed out as you do.

The second part of the answer is that we strongly encourage you to initialize as many of your data members as possible using the colon initializer syntax. This is simply a C++ best practice: if at all possible, do not allow uninitialized or incompletely initialized variables of any kind.

**Listing 12.1:** The output produced by Optional\_module.cc when run using optional.fcl

```
1
2 Hello from Optional::constructor.
3 Hello from Optional::beginJob.
4 Hello from Optional::beginRun: run: 3
5 Hello from Optional::beginSubRun: run: 3 subRun: 0
6 Hello from Optional::analyze. Event id: run: 3 subRun: 0 event: 1
7 Hello from Optional::analyze. Event id: run: 3 subRun: 0 event: 2
8 Hello from Optional::analyze. Event id: run: 3 subRun: 0 event: 3
9 Hello from Optional::analyze. Event id: run: 3 subRun: 0 event: 4
10 Hello from Optional::analyze. Event id: run: 3 subRun: 0 event: 5
11 Hello from Optional::beginSubRun: run: 3 subRun: 1
12 Hello from Optional::analyze. Event id: run: 3 subRun: 1 event: 1
13 Hello from Optional::analyze. Event id: run: 3 subRun: 1 event: 2
14 Hello from Optional::analyze. Event id: run: 3 subRun: 1 event: 3
15 Hello from Optional::analyze. Event id: run: 3 subRun: 1 event: 4
16 Hello from Optional::analyze. Event id: run: 3 subRun: 1 event: 5
17 Hello from Optional::beginSubRun: run: 3 subRun: 2
18 Hello from Optional::analyze. Event id: run: 3 subRun: 2 event: 1
19 Hello from Optional::analyze. Event id: run: 3 subRun: 2 event: 2
20 Hello from Optional::analyze. Event id: run: 3 subRun: 2 event: 3
21 Hello from Optional::analyze. Event id: run: 3 subRun: 2 event: 4
22 Hello from Optional::analyze. Event id: run: 3 subRun: 2 event: 5
```

One reasonable guideline is that physics related tasks belong in the beginJob member function while computer science related tasks belong in the constructor. Your experiment may have additional guidelines.

For those of you familiar with ROOT, we can provide an example of something physics-related. We suggest that you create histograms, ntuples or trees in one of the begin member functions, perhaps beginJob or beginRun. Other constraints may enter into this decision. If booking a histogram requires geometry information to set the limits correctly, that information may be run-dependent and you will need to book these histograms in beginRun, not beginJob.



## 12.9 Suggested Activities

## 12.9.1 Add the Matching end Member functions

art defines the following three member functions:

```
void endJob () override;
void endRun ( art::Run const& run ) override;
void endSubRun ( art::SubRun const& subRun ) override;
```

Go to your source window. In the file Optional\_module.cc, add these member functions to the declaration of the class Optional and provide an implementation for each. In your implementation, just copy the printout created in the corresponding begin function and, in that printout, change the string "begin" to "end".

Then go to your build window and make sure that your current directory is your build directory. Then rebuild this module and run it:

```
buildtool art -c fcl/OptionalMethods/optional.fcl >& output/optional2.log
```

Consult Chapter 10 if you need to remember how to indentify that the build completed successfully. Compare the output from this run of *art* with that of the previous run: do you see the additional printout from the member functions that you added?

The solution to this activity is provided as the file OptionalAnswer01\_module.cc. It is already built. You can run it with:

## art -c fcl/OptionalMethods/optionalAnswer01.fcl >& output/optionalAnswer01.log

Does the output of your code match the output from this code?

### 12.9.2 Run on Multiple Input Files

In a single run of *art*, run your modified version of the module Optional on all of the three of the following input files:

```
inputFiles/input01_data.root
inputFiles/input02_data.root
inputFiles/input03_data.root
```

If you need a reminder about how to tell *art* to run on three input files in one job, consult Section 9.8.5.

Make sure that the printout from this job matches the description of the event loop found in Section 3.6.2.

## 12.9.3 The Option –trace

The *art* command supports a command line option named -trace. This creates additional printout that identifies every step in the event loop. Use this option to trace what *art* is doing when you run this exercise. For example

```
art -c fcl/OptionalMethods/optional.fcl -trace >& output/trace.log
```

You should be able to identify your printout among the printout from *art* and see that your printout appears in the expected place.

When you are getting an error from *art* and you don't understand which module is causing the problem, you can use -trace to narrow your search.

# 13 Exercise 4: A First Look at Parameter Sets

## 13.1 Introduction

In the previous few chapters you have used FHiCL files to configure art jobs. Recall from Section 9.8 that in a FHiCL file (a group of FHiCL definitions delimited by braces  $\{\ \}$ ) is called a table. When art reads its run time configuration FHiCL file, it transforms the FHiCL file into a C++ representation; in that representation, a FHiCL table becomes an object of type fhicl::ParameterSet, which we refer to as a  $parameterset(\gamma)$ .

Among other things, you have learned how to define a module label and its corresponding parameter set, the simplest case looking like:

```
moduleLabel : {
   module_type : ClassName
}
```

where the moduleLabel is an identifier that you define and ClassName is the name of a module class. *art* requires that the module\_type parameter be present.

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When you define a module label, you may enter additional FHiCL definitions (i.e., *parameters*) between the braces to form a larger parameter set. For example:

This functionality allows you to write modules whose behaviour is run-time configurable. For example, if you have a reconstruction algorithm that depends on some cuts, the values of those cuts can be provided in this way.

## 13.2 Prerequisites

The prerequisite for this chapter is all of the material in Part I (Introduction) and the material in Part II (Workbook) up to and including Chapter 10. You can read this chapter without necessarily having read Chapter 11 or 12.

## 13.3 What You Will Learn

In Section 10.7.3.4 you saw that the constructor of a module is required to take an argument of type fhicl::ParameterSet const&.

In this chapter you will learn how to use this argument to read additional parameters in a parameter set. In particular, you will learn about the class fhicl::ParameterSet and after working through the exercises in this section, you should know how to:

1. read parameter values from a FHiCL file into a module

- 2. require that a particular parameter be present in a parameter set
- 3. use data members to communicate information from the constructor to other member functions of a module
- 4. print a parameter set
- 5. use the colon initializer syntax
- 6. provide a default value for a parameter (if the parameter is absent from a parameter set)
- 7. modify the precision of the printout of floating point types
- 8. recognize the error messages for a missing parameter or for a value that cannot be converted to the requested type

#### You will also learn:

- 1. that you should find out your experiment's policy about what sorts of parameters are allowed to have default values
- 2. the two special parameters automatically provided by art
- 3. the canonical forms of parameters

Finall, you will learn a small amount about C++ templates and C++ exceptions, just enough to understand the exercise.

## 13.4 Setting up to Run this Exercise

To run this exercise, you need to be logged in to the computer on which you ran Exercise 2 (in Chapter 10). If you are continuing on from a previous exercise, you need to keep both your source and build windows open.

If you are logging back in, follow the instructions in Section 10.5 to reestablish your source and build windows.

In your source window, cd to your source directory. Then cd to the directory for this exercise and look at its contents:

cd art-workbook/ParameterSets

ls

```
CMakeLists.txt pset02.fcl PSet03_module.cc
pset01.fcl PSet02_module.cc pset04.fcl
PSet01_module.cc pset03.fcl PSet04_module.cc
```

The source code for the first module you will run is PSet01\_module.cc and the FHiCL file to run it is pset01.fcl. The file CMakeLists.txt is identical to that used by the previous two exericses. The remaining files are the source and FHiCL files for additional steps in this exercise.

In your build window, make sure that you are in your build directory. At this time you do not need to build any code because all code for the Workbook was built the first time that your ran buildtool.

## **13.5** The Configuration File pset01.fcl

The FHiCL file that you will run in this exericse is pset01.fcl. Look at this file in your source window. You will see that pset01.fcl defines a parameter set psetTester, shown below, to configure an analyzer module named PSet01.

```
analyzers: {
    psetTester : {
        module_type : PSet01
        a : "this is quoted string"
        b : 42
        c : 3.14159
        d : true
        e : [ 1, 2, 3 ]
        f : {
            a : 4
            b : 5
        }
    }
}
```

Figure 13.1: The FHiCL definition of the parameter set psetTester from pset01.fcl.

This parameter set tells *art* to run the module PSet01; it then supplies some additional parameter definitions that will be read by that module. Additional definitions like these in a FHiCL file have the following properties:

- 1. The module specified by the module\_type parameter defines which parameters must be present in this list, and which parameters are optional.
- 2. Each definition must be a legal FHiCL definition.
- 3. These definitions have no meaning, per se, to *art* or to FHiCL; they are used by the C++ code.
- 4. Each definition may use the full power of FHiCL and my contain nested parameter sets to arbitrary depth.

Looking at the parameter set, it appears that the parameter a has a value that is a string of text, parameter b's value is an integer number, parameter c's is a floating point number, parameter d's is one of the two possible boolean values, parameter e's is an array of integers and that parameter f's is a nested parameter set. You will learn in Section 13.6 that the (eventual) instantiation of the class PSet01 internally represents the value of

each parameter a a string. The computer-science-speak for this is that FHiCL is a *type-free* language.

## 13.6 The Source code file PSet01\_module.cc

The source code for this exercise is found in the file PSet01\_module.cc. The new features seen in this exercise are all in the definition of the constructor. The FHiCL file will tell *art* to call this module and will set the variable pset (represented by this argument) to the values in the parameter set psetTester.

The first change is that the constructor now *uses* its argument (i.e., refers to it within the declaration); therefore the argument must have a name, in this case pset. Any name could have been chosen for this argument. Let's examine the first part of the constructor.

The type (class) of the argument is fhicl::ParameterSet const&. This class ParameterSet is in the namespace fhicl; all identifiers in the namespace fhicl are found in the UPS package named **fhiclcpp**. You can find the header file for ParameterSet at \$FHICLCPP\_INC/fhiclcpp/ParameterSet.h.

The argument named pset is passed by const reference; i.e., the module is not allowed to modify it (due to const), and the module is given access to it, but not given a copy of it (due to the &, meaning "reference").

Parameters are set below, with the name of the parameter and the type of variable that will receive the returned value. The syntax used to specify the return type uses a feature of C++ called  $templates(\gamma)$ . This section of code is further discussed within the chapter text.

```
std::string a = pset.get<std::string>("a");

int b = pset.get<int> ("b");

double c = pset.get<double>("c");

bool d = pset.get<bool> ("d");

std::vector<int> e = pset.get< std::vector<int> >("e");

fhicl::ParameterSet f = pset.get<fhicl::ParameterSet>("f");

int fa = f.get<int>("a");

int fb = f.get<int>("b");
```

After copying psetTester into pset, art adds a first parameter module\_type (below) and sets it to the identifier module\_type, and adds a second parameter module\_label that it sets it to the module label as defined in the FHiCL file, in this case psetTester. This extra step is not done for every parameter set defined within a FHiCL file; it is only done for parameter sets that are used to configure modules.

Recall from Section 13.5 that the instantiated object PSet01 internally represents the value of each parameter as a string. If you ask that the value of a parameter be returned as a string, pset will simply return a copy of its internal representation of that parameter. On the other hand, if you ask that the value of a parameter be returned as any other type, then pset needs to do some additional work. For example, if you ask that a parameter be returned as an int, then pset must first find its internal string representation of that parameter; it must then convert that string into a temporary variable of the requested type and return the temporary variable. Therefore, when your code asks the variable pset to return the value of a parameter, it must tell pset two things:

- 1. the name of the parameter
- 2. the type of the variable that will receive the returned value

The angle bracket syntax <> used in this example to specify the return type is the signature of a feature of called  $templates(\gamma)$ . art and FHiCL use templates in several prominent places. The name of the parameter is specified as a familiar function argument while the return type is specified between the angle brackets. The name between the angle brackets is called a template argument. If you do not supply a template argument, then your code will not compile.

You do not need to fully understand templates — just how to use them when you encounter them. For example, the line that sets the parameter a that reads

first declares a local variable named a that is of type std::string and then does the following:

- 1. It asks pset if it has a parameter named a.
- 2. If it has this parameter, pset will return it as a string.
- 3. The returned value is used to initialize the local variable named a.

Section 13.9 will describe what happens if pset does not have a parameter named a.



It is not required that the local variable, a, have the same name as the FHiCL parameter a. But, with rare exceptions, it is a good practice to make them either exactly the same or, at a minimum, obviously related.

The following line, that sets the parameter b, that reads

```
int b = pset.get<int>("b");
```

is similar to the previous line; the main difference is that pset will convert the string to an int before returning it. pset knows that it must perform the conversion to int because the template argument tells it to.

It is beyond the scope of this chapter to discuss how the template mechanism is used to trigger automatic type conversions. It is sufficient to always remember the following: when you use the get member function of the class fhicl::ParameterSet, the template argument must always match the type of the variable on the left hand side. Templates will be discussed in Section 13.8.



The authors of FHiCL could have designed a different interface, such as

```
std::string a = pset.get_as_string("a");
std::string b = pset.get as int ("b");
```

Instead they chose to write it using templates. The reason for this choice is that it allows one to add new types to FHiCL without needing to recompile FHiCL. How you do this is beyond the scope of this chapter. You now know everything that you need to know about templates in order to use fhicl::ParameterSet effectively.

The rest of the lines in that section of code extract the remaining parameters from pset and make copies of them in local variables. The remainder of the constructor, shown below, prints the values of these parameters; these make the printout in lines 1 through 11 of Listing 13.1.

```
std::cout « "\n----\nPart 1:\n";
std::cout « "a : " « a « std::endl;
std::cout « "b : " « b « std::endl;
std::cout « "c : " « c « std::endl;
std::cout « "d : " « d « std::endl;
std::cout « "e :":
for ( int i: e ){
  std::cout « " " « i; }
std::cout « std::endl;
    The following lines show one way to print the two values,
    a and b from the parameter set f.
std::cout « "f.a : " « fa « std::endl:
std::cout « "f.b : " « fb « std::endl;
std::cout « "module_type: " « module_type « std::endl;
std::cout « "module_label: " « module_label « std::endl;
    The following three lines show show two other ways, using the to_string()
    and to indented string() member functions of the class
    fhicl::ParameterSet. These lines make the printout found in lines 13
    through 18 of Listing 13.1.
std::cout « "\n——\nPart 2:\n";
std::cout « "f as string: " « f.to_string() « std::endl;
std::cout « "f as indented-string:\n" « f.to_indented_string() « std::endl;
std::cout « "\n———\nPart 3:\n";
std::cout « "pset:\n" « pset.to_indented_string() « std::endl;
    The last two lines use the to indented string() member function to print
    everything found in the parameter set psetTester. These lines make the printout
    found in lines 21 through 36 of Listing 13.1.
```

Your code may ask for the values of parameters from a ParameterSet in any order, and any number of times, including zero.

One final comment on PSet01 module.cc: the analyze member function is empty.

Nevertheless, it must be present because *art* requires all analyzer modules to provide a member function named analyze. If we removed this member function from the class PSet01, then the module would not compile.

# 13.7 Running the Exercise

Now let's see what happens when you run the job. In your build directory, run the following command

```
art -c fcl/ParameterSets/pset01.fcl >& output/pset01.log
```

The expected output from this command is shown in Listing 13.1.

The module reads in the parameter set and then prints out each of the values in several different ways. Check that the printout matches the definitions of the parameters from pset01.fcl.

# 13.8 Member Function Templates and their Arguments

Now that you have seen templates, we can introduce some more language that you will need to know. In the above examples, get<std::string> and get<int> are member functions of the class ParameterSet (represented in the module by the argument pset).

On its own, get is called a *member function template*; this means that get is a set of rules to write a member function. The member function can only be written once the template's argument has been specified. In the future, when we refer to get, we will call it by its proper name:

```
ParameterSet::qet<T>
```

or, sometimes, just get<T>. In the notation <T>, the angle brackets indicate that get is a template and the capital letter T is a dummy argument that indicates that if you want to use the template, you must supply one template argument. The choice of the letter T as the name of the dummy argument is a mnemonic for Type, indicating that the template argument must be the name of a type.

**Listing 13.1:** Output from PSet 01 when run using pset 01.fcl (without extraneous messages)

```
1 -----
2 Part 1:
3 a : this is quoted string
4 b: 42
5 c: 3.14159
6 d:1
7 e: 1 2 3
8 f.a: 4
9 f.b: 5
10 module_type: PSet01
11 module_label: psetTester
12
13 -----
14 Part 2:
15 f as string: a:4 b:5
16 f as indented-string:
17 a: 4
18 b: 5
19
20
21 -----
22 Part 3:
23 pset:
24 a: "this is quoted string"
25 b: 42
26 c: 3.14159
27 d: true
28 e: [ 1
29
30
31
]
32 f: { a: 4
33 b: 5
34 }
35 module_label: "psetTester"
36 module_type: "PSet01"
```

# 13.9 Exceptions

#### 13.9.1 Error Conditions

There are two sorts of error conditions that may occur when reading parameters from a parameter set:

- 1. The requested parameter is not present in the parameter set.
- 2. The requested parameter is present but cannot be converted into the requested type.

To give an example of the second sort, suppose that on line 6 of Listing ?? someone changed the FHiCL definition of the parameter c from 3.14159 to "test". Now consider what happens when line 6 of Listing ?? is executed: the code will correctly find that parameter c exists but it will produce an error when it tries to convert the string "test" to a double.

#### 13.9.2 Error Handling

The C++ language has a feature called  $exceptions(\gamma)$ . As for templates, this is an advanced feature that you do not need to understand in detail but you do need to be aware of it in broad strokes. In this case, exceptions are used behind the scenes and you will not see any code that explicitly uses exceptions.

When an error of this sort occurs, *art*'s default response is to abort the processing of your module and to return control to *art*, which will then attempt an orderly shutdown, meaning that it will:

- 1. Record that your module had an error condition that stopped processing; this information will be written to all output event-data files.
- 2. Write a message to the log file that describes what happened.
- 3. Call the endSubRun member function of every module.
- 4. Call the endRun member function of every module.
- 5. Call the endJob member function of every module.
- 6. Properly flush and close all output and log files.

- 7. Perform a few other clean up and shutdown actions for parts of *art* that have not yet been discussed.
- 8. Return a non-zero status code to the parent process; the status code is the number that appears on the last line of your *art* output, the line that begins "Art has completed ...".

For most sorts of errors, the orderly shutdown will be successful and your work up to the exception will be preserved.

But there are circumstances for which the orderly shutdown will fail. One example of this is if you have reached your disk quota and there is no disk space to hold more output.



art provides a mechanism that allows the user to specify, at run time, alternate behavior. For example, there may be some errors for which you wish to write the offending event to a separate output file and to continue with the processing of the next event. For some other errors you may wish that art continue the current event and procede to the next module. If you do not specify an alternate behavior, art will perform its default action, which is to attempt an orderly shutdown. A detailed discussion of these features is beyond the scope of this chapter.

#### 13.9.3 Suggested Exercises

In pset01.fcl, remove the definition of the parameter b. Rerun *art*. You should see an error message like that shown in Listing 13.2. Read the error message and understand what it is telling you so that you will recognize the error message if you make this mistake in the future.

**Listing 13.2:** The expected output when the parameter b is removed from pset01.fcl

Part

In pset01.fcl, restore the definition of b and change the definition of c to "test". Rerun *art*. You should see an error message like that shown in Listing 13.3. Again, read the error message and understand it.

**Listing 13.3:** The expected output when the parameter c misdefined in pset 01.fcl

```
%MSG-s ArtException: PSet01:psetTester@Construction 14-Jul-2013 19:42:54 CDT
   ModuleConstruction
2 cet::exception caught in art
3 ---- Type mismatch BEGIN
5
    ---- Type mismatch BEGIN
6
      error in float string:
7
      test
8
      at or before:
9
    ---- Type mismatch END
10 --- Type mismatch END
11 %MSG
12 Art has completed and will exit with status 8001.
```

#### 13.10 Parameters and Data Members

Very often information from the parameter set is needed in a member function of the module class. The way to propagate this information from the parameter set to the member function is to store the values of these parameters as data members of the module class. This is illustrated in the two files PSet02\_module.cc and pset02.fcl.

Refer back to Section 6.6.5 to refamiliarize yourself with data members or with the colon initializer syntax, as needed. Notice the three (private) member functions listed with their return types, and that they are used in the analyze block.

To run this example, enter

```
\mathsf{art}\, \mathsf{-c}\, \mathsf{fcl}/\mathsf{ParameterSets}/\mathsf{pset02}.\mathsf{fcl} > \& \, \mathsf{output/pset02}.\mathsf{log}
```

The expected output from this is given in Listing 13.4.

**Listing 13.4:** The output from PSet 02 when run using pset 02.fcl

```
1 Event number: run: 1 subRun: 0 event: 1 b: 42 c: 3.14159 f: a:4 b:5
2 Event number: run: 1 subRun: 0 event: 2 b: 42 c: 3.14159 f: a:4 b:5
3 Event number: run: 1 subRun: 0 event: 3 b: 42 c: 3.14159 f: a:4 b:5
```

This example is only relevant when parameters are actually used in member functions. If a parameter is used only inside the constructor, do not store it as a data member; instead you should store it as a local variable of the constructor. This brings up a "best practice:" always declare a variable in the narrowest scope that works.



# 13.11 Optional Parameters with Default Values

It is sometimes convenient to provide a default value for a parameter. Default values may be provided in the source code that reads the parameter set. This mechanism is illustrated by the files PSet03\_module.cc and pset03.fcl.

You have already seen that the member function template ParameterSet::get<T> takes one function argument, the name of the parameter. For example,

```
int b = pset.get<int>("b");
```

It also takes an optional second function argument, a default value for the parameter. For example,

```
int b = pset.get < int > ("b", 0);
```

If the second argument is present, there two cases:

- 1. If the parameter is not defined in pset , then the second argument is returned as the value of the call to get.
- 2. If the parameter is defined in pset, then the second argument is ignored and the value read from the FHiCL file is returned as the value of the call to get.

When reading the code in this example you will encounter the expression:

```
std::vector<double>(5,1.0)
```

This tells the compiler to instantiate an object of type std::vector<double>, set its size to 5 and initialize elements 0 through 4 to have the value 1.0. If you are not familiar with this syntax, you can read about it in the STL documentation (see Section 6.7).

This expression appears as the second argument of a call to pset.get. Therefore the compiler will create an unnamed temporary object (the vector of doubles) and pass that

**Listing 13.5:** The parameter-related portion of the output from PSet03 when run using pset03.fcl

```
1 debug level: 0
2 g: 1 1 1 1 1
```

object as the second argument; once the function call has completed, the temporary object will be deleted.

With the above explanations, the source code for this example should be reasonably self-explanatory; it looks for two parameters named debugLevel and g and supplies default values for each of them. Look at the file pset03.fcl; you will see that the parameters debugLevel and g are not present in the testPSet parameter set.

To run this example,

```
art -c fcl/ParameterSets/pset03.fcl >& output/pset03.log
```

The expected output from this is given in Listing 13.5.

As a suggested exercise, edit pset03.fcl and, in the parameter set testPSet, provide definitions for the parameters debugLevel and g. Make their values different from the default values. Rerun *art* and verify that the module has correctly read in and printed out the values you defined.

#### 13.11.1 Policies About Optional Parameters

Allowing *optional* parameters is important for developing, debugging and testing; if all parameters were required all of the time, the complete list of parameters could become unwieldy\*. On the other hand, the use of optional parameters can make it difficult to audit the physics content of a job. Therefore experiments typically have policies for what sorts of parameters may have defaults and what sorts may not. For example, your experiment may prohibit default values for parameters that define the physics behavior, but allow them for other parameters.



Consult your experiment to learn what policies you should follow.

<sup>\*</sup>FHiCL has several features that make it easier to deal with large parameter sets. *This will be explained in a future chapter.* 

# 13.12 Numerical Types, Precision and Canonical Forms

FHiCL recognizes numbers in both fixed point and exponential notation, for example 123.4 and 1.234e2; the letter e that separates the exponent can be written in either upper or lower case.

In the preceding exercises you defined some numerical values in a FHiCL file, read them into your code and printed them out; the printed values exactly matched the input values. The values used in those exercises were carefully chosen to avoid a few surprises: there are cases in which the printed value will be an equivalent, but not identical, form. This section discusses those cases and provides some examples.

When FHiCL recognizes that a parameter value is a number it converts the number into a *canonical form* and stores the canonical form as a string. The transformation to the canonical form preserves the full precision of the number and involves the following steps:

- 1. The canonical form has no insignificant characters:
  - (a) no insignificant trailing zeros
  - (b) no insignificant trailing decimal point
  - (c) no insignficant leading plus sign
  - (d) no insignficant leading plus sign in the exponent
- 2. If a number is specified in exponential notation and if the number can be represented as a integer without loss of precision, and if the resulting integer has 6 or fewer digits, then the canonical form is the integer. For example, the canonical form of 1.23456E5 is 123456 but the canonical form of 1.23456E6 is 1.23456e6.
- 3. The canonical form of all other floating point numbers is exponential notation with a single, non-zero digit to the left of the decimal point.
- 4. The canonical form of all strings includes beginning and ending quotes; this is true even if the string contains no embedded whitespace or other special characters.

Some examples of numbers and their canonical forms are given in Table 13.1.

If a numerical value, when expressed as a fixed point number, has no fractional part, your code may ask for the parameter to be returned as either a floating point type (

Number Canonical Number Canonoical Form Form 2 2 1.234E2 1.234e2 2. 2 123456 1.23456E5 2.0 2 1.23456E6 1.23456e6 2.1E2 210 1234567 1.234567e6 +210 210 0.01 1E-2

Table 13.1: caption

such as double or float ) or as an integral type (such as int, short unsigned or std::size\_t). For example, line 6 of Listing ?? could have been written:

```
double b = pset.get<double>("b");
```

This code will do the expected thing: in the example pset01.fcl, it will read the value 42 into a variable of type double.

On the other hand, if a numercial value, when expressed as a fixed point number, does have a fractional part, you may only ask for the parameter to be returned as a floating point type. If you ask for such a value as an int, the ParameterSet::get<int> member function will throw an exception; similarly for all other integral types. This behavior may not be intuitive: the authors of *art* could have decided, instead, to discard the fractional part and return the integer part. They chose not to do this because when this situations occurs, it is almost always an error.

#### 13.12.1 Suggested Exercises

The above ideas are illustrated by the files PSet04\_module.cc and pset04.fcl. To run this example,

```
art -c fcl/ParameterSets/pset04.fcl >& output/pset04.log
```

The expected output from this is given in Listing 13.6. Read the source code and the FHiCL file; then examine the output. The first three lines show three different printed formats of the parameter a, with the first being the canonical form. While all forms are equal to the number found in the FHiCL file, they all have different formats. Understand why each line has the format it does.

**Listing 13.6:** The output from PSet04 when run using pset04.fcl

```
1 parameter a as a string: 1.23456e6
2 parameter a as a double: 1.23456e+06
3 parameter a as int: 1234560
4 parameter b as a string: 3.1415926
5 parameter b as a double: 3.14159
6 parameter b as a double with more significant figures: 3.1415926
7 parameter c as a string: 1
8 parameter c as an int: 1
```

Line 4 shows the canonical form of the parameter b. Line 5 shows what is printed using the default C++ settings; the two least significant characters were dropped. The code that produces line 6 shows how the use the STL precision function to tell C++ to print more significant figures.

If you modify the precision of cout, it will change the format of the printout for the rest of the job; usually this is a bad thing. To avoid this, PSet04\_module.cc illustrates how to save and restore the precision of cout.

Line 7 shows the canonical form of the parameter c and line 8 shows the default C++ printed form of the integer.

For the next exercise, edit pset04.fcl and change the value of c to something with a fractional part. Rerun *art*; you should see that it throws an exception because it is illegal to read a numeric value with a fractional part into a variable of integral type. The error message from *art* is shown in Listing 13.7. Read the error message and understand what it is telling you so that you will recognize the error message if you make this mistake yourself.

**Listing 13.7:** The output from PSet04 when run using the modified version of pset04.fcl, which has an intentional error

# 14 Exercise 5: Making Multiple Instances of a Module

#### 14.1 Introduction

In a typical HEP experiment is often necessary to repeat one analysis several times, with each version differing only in the values of some cuts; this is frequently done to tune cuts or to study systematic errors. Very often it is both convenient and efficient to run all of the variants of the analysis in a single job.

A powerful feature of *art* is that it permits you to run an *art* job in which you define and run many instances of the same module; when you do this, each instance of the module gets its own parameter set. In this chapter you will learn how to use this feature of *art*.

## 14.2 Prerequisites

The prerequisite for this chapter is all of the material in Part I (Introduction) and the material in Part II (Workbook) up to and including Chapter 13, but excluding Chapter 12.

### 14.3 What You Will Learn

In this chapter you will learn how to run an *art* job in which you run the same module more than once. This exercise will make it clear why *art* needs to distinguish the two ideas of *module label* and module\_type.

## 14.4 Setting up to Run this Exercise

To run this exercise, you need to be logged in to the computer on which you ran Exercise 2 (in Chapter 10). If you are continuing on from a previous exercise, you need to keep both your source and build windows open.

If you are logging back in, follow the instructions in Section 10.5 to reestablish your source and build windows.

In your source window, cd to your source directory. Then cd to the directory for this exercise and look at its contents

```
$ cd art-workbook/ModuleInstances
$ ls
CMakeLists.txt magic.fcl MagicNumber_module.cc
```

The source code for the first module you will run is MagicNumber\_module.cc and the FHiCL file to run it is magic.fcl. The file CMakeLists.txt is identical that used by the previous two exericses.

In your build window, make sure that you are in your build directory. At this time you do not need to build any code becaue all code for the Workbook was built the first time that your ran buildtool.

## 14.5 The Source File Magic\_module.cc

The source code for this exercise is found in the file Magic\_module.cc. Look at this file and you should see the following features, all of which you have seen before.

- 1. The file declares and defines a class named MagicNumber that follows the rules to be an *art* analyer module.
- 2. The class has a constructor and an analyze method.
- 3. The class has a data member named magicNumber\_, of type int.
- 4. The class initializes magicNumber\_by reading a value from its parameter set; the name of the parameter is magicNumber (without the underscore).

- 5. The parameter magic Number is a required parameter.
- 6. Both the constructor and the analyze method print an informational message that includes the value of magicNumber\_.

## 14.6 The FHiCL File magic.fcl

The FHiCL file used to run this exercise is magic.fcl. Look at this file and you should see the following features:

- 1. Compared to previous exercises, The FHiCL names process\_name, source and services have no important differences.
- 2. In the analyzers parameter set, inside the physics parameter set, you will see the definition of four module labels, boomboom, rocket, flower and bigbird\*. The value of each definition is a parameter set.
- 3. The first three of these parameter sets tell *art* to run the module MagicNumber and each provides a value for the required magicNumber parameter<sup>†</sup>
- 4. The last parameter set tells *art* to run the module First, the source for which was discussed in Chapter 10; this module does not need any additional parameters.
- 5. The path el contains the names of all of the module labels from the analyzers parameter set.

## 14.7 Running the Exercise

In your build directory, run the following command

```
$ art -c fcl/ModuleInstances/magic.fcl >& output/magic.log
```

The expected output from this command is shown in Listing 14.1; for clarity, the printout made by *art* has been elided. Compare this printout to the printout from your run; it should

<sup>\*</sup> All of these are nicknames of ice hockey players who played for the Montreal Canadiens ice hockey team; all of them have had their sweater number retired

<sup>†</sup>In each case the magic number is the sweater number of the hockey player whose nickname is the module label.

Listing 14.1: The output from running magic.fcl

```
1 MagicNumber::constructor: magic number: 9
2 MagicNumber::constructor:
                             magic number: 5
3 Hello from First::constructor.
4 MagicNumber::constructor: magic number: 10
5 MagicNumber::analyze: event: run: 1 subRun: 0 event: 1 magic number: 9
6 MagicNumber::analyze: event: run: 1 subRun: 0 event: 1 magic number: 5
7
  Hello from First::analyze. Event id: run: 1 subRun: 0 event: 1
8 MagicNumber::analyze: event: run: 1 subRun: 0 event: 1 magic number: 10
9 MagicNumber::analyze: event: run: 1 subRun: 0 event: 2 magic number: 9
10 MagicNumber::analyze: event: run: 1 subRun: 0 event: 2 magic number: 5
11 Hello from First::analyze. Event id: run: 1 subRun: 0 event: 2
12 MagicNumber::analyze: event: run: 1 subRun: 0 event: 2 magic number: 10
13 MagicNumber::analyze: event: run: 1 subRun: 0 event: 3 magic number: 9
14 MagicNumber::analyze: event: run: 1 subRun: 0 event: 3 magic number: 5
15 Hello from First::analyze. Event id: run: 1 subRun: 0 event: 3
```

be exactly the same. Inspect the printout and the files MagicNumber\_module.cc and ../FirstModule/First\_module.cc; understand why the printout is what it is.

#### 14.8 Discussion

#### 14.8.1 Order of Analyzer Modules is not Important

As it happens, *art* runs the four analyzer modules in the order specified in the path definition e1. But you must not count on this behaviour! Two of the design rules of *art* are:



- 1. Modules may only communicate with each other by putting information into, and reading information from, the art::Event.
- 2. Analyer modules may not put information into the art::Event.

Therefore *art* is free to run analyzer modules in any order.

For producer modules, which may add information to the event, the order of execution is often very important. When you reach the exercises that run producer modules, you will be told how to specify the order of execution.

You may wish to review some of the other ideas about *art* paths that are described in Section 9.8.8.

#### 14.8.2 Two Meanings of Module Label

In the preceding discussion, the name *module label* was used in two subtlely different ways, as is illustrated by the module label rocket:

- rocket identifies a parameter set that is used to configure an instance of the module MagicNumber.
- 2. rocket is also used as the name of the module instance that is configured using this parameter set; the elements in the path e1 are all the names of module instances.

Clearly these two meanings are very closely related, which is why the same name, *module label*, is used for both ideas. Throughout the remainder of this document suite the name *module label* will be used for both meanings; the authors believe it will be clear from the context which meaning is intended. This is standard usage within the *art* community.

# 14.9 Suggested Exercise

Edit magic.fcl and do the following:

- 1. Add a new analyzer module label that configures an instance of the module Optional from Chapter 12.
- 2. Add the new module label to e1.

Then re-run magic.fcl. Do you see the expected additional printout?

#### 14.10 Review

After working through this exercise, you should:

- 1. Know how to run multiple instances of the same module within one art job.
- 2. Understand that *art* does not guarantee the order in which analyzer modules will be run.

3. Understand the two senses in which the name *module label* is used: as the name of a parameter set and as the name of the corresponding instance of a module.

# 15 Exercise 6: Accessing Data Products

#### 15.1 Introduction

Section 10.7.3.6 described the class art::Event as an art::EventID plus a collection of data products. The concept of a data product was described in Section 3.6.4. You have already done several exercises that made use of the art::EventID and in this chapter you will do your first exercises that use a data product.

# 15.2 Prerequisites

Prerequisites for this chapter include all of the material in Part I (Introduction) and the material in Part II (Workbook) up to and including Chapter 13.

You must also be familiar with the toy experiment described in Section 3.7.

This exercise will use class templates and member function templates in several places. The use of templates was introduced in Section 13.6. You need to know how to use templates but you do not need to know how to write one. You will need a minimal understanding of the class template std::vector, which is part of the C++ Standard Library. If you understand the following four points, then you understand enough about std::vector for this exercise. If t is an object of type std::vector<T>, then:

- 1. t behaves much like an array of objects of type T. The main difference is that capacity of the array automatically grows to be large enough to hold all of the elements in the array.
- 2. The identifier inside the angle brackets is called a *template argument* and it is usually

the name of a C++ type. \*

- 3. The dynamic sizing occurs in the middle of a running program; not at compile time.
- 4. This expression sets nEntries to the number of entries in t:

```
std::size_t nEntries = t.size();
```

15.3 What You Will Learn

In this exercise you will learn about:

```
1. the data type tex::GenParticleCollection
```

- 2. the four-part name of an art data product
- 3. the class art::InputTag
- 4. the class template art::Handle
- 5. the class template art::ValidHandle
- 6. the member function templates of art::Event:

```
o getByLabel( art::InputTag, art::Handle<T>) const;
o getValidHandle<T>( art::InputTag ) const;
```

# 15.4 Background Information for this Exercise

The input files used for the *art* workbook contain data products created by a workflow that simulates the response of the toy detector to a generated event, described in Section 3.7.2. The first step in this workflow is to create the generated event, which is stored in the art::Event as data product.

<sup>\*</sup>You will only see one case in the entire workbook in which it is something other than a the name of a C++ type; and this will be during a short side trip to discuss the **cetlib** utility library.

Listing 15.1: The contents of GenParticleCollection.h

```
#include "toyExperiment/MCDataProducts/GenParticle.h"

#include <vector>

namespace tex {

typedef std::vector<GenParticle> GenParticleCollection;
}
```

In this exercise you will retrieve this data product from the art::Event and print the number of generated particles in each event. In Chapter 17, you will look at the properties of individual generated particles.

#### **15.4.1 The Data Type** GenParticleCollection

Each generated particle in the simulated event is described by an object of type tex::GenParticle. All of the generated particles in a given event are stored in an object of type tex::GenParticleCollection. This object is written to the art::Event as a data product.

The header files that describe these two classes, GenParticle.h and GenParticleCollection.h, are found under:

```
$TOYEXPERIMENT_INC/toyExperiment/MCDataProducts/
```

The content of GenParticleCollection.h is shown in Listing 15.1; the code guards and comments have been omitted. This header uses a typedef to declare that the name tex::GenParticleCollection; is a synonym for std::vector<tex::GenParticleCollection>.

Why did the authors of the workbook decide to use a typedef and not simply ask you to code std::vector<GenParticle> when needed? The reason is future-proofing. Suppose that down the road the authors find that they need to change the definition of tex::GenParticleCollection; if you used the typedef, it is much more likely

that your code will continue to compile and work correctly as is. If, on the other hand, you used std::vector<GenParticle>, then you would need to identify and edit every instance.

Please use the typedef GenParticleCollection in your own code and do not hand-substitute its definition.



did authors of the workbook decide Why the to call this typedef GenParticleCollection and not, for example, GenParticleVector? The answer is a different sort of future-proofing. The C++ standard library provides class templates other than vectors that are collections of objects, and one can imagine a scenario in which it would make sense to change GenParticleCollection to use a collection type, such as std::deque for example. In such a scenario, the following definition would make perfect sense to the C++ compiler but would be misleading to human readers:

typedef std::deque<GenParticle> GenParticleVector The generic name *Collection* avoids this problem.

#### 15.4.2 Data Product Names

Each *art* data product has a name that is a text string with four fields, delimited by underscore characters (\_) that represent, in order, the data type, module label, instance name and process name, e.g.,:

MyDataType\_MyModuleLabel\_MyInstanceName\_MyProcessName

Each data product name must be unique within an *art* event-data file. The fields in the data product name may only contain the following characters $^{\dagger}$ :

- a...zA...z
- 0 0 . . . 9
- :: (double colon)

<sup>†</sup>Experts may want to know that in an *art* event-data file, each data product is stored as a TBranch of a TTree named Event. Only these characters are legal in a TBranch name. The name of the TBranch is the name of the data product, hence the restriction.

In particular, periods, dashes, commas, underscores, semicolons, white space and single colons are not allowed; underscores are only allowed as the field separator, not within a field.

#### About each field:

- 1. The data type field is the so-called *friendly name* name of the data type for the data product; friendly names are discussed below.
- 2. The module label field is the label of the module that created the data product. Note that it is the module *label* as specified in the FHiCL file, not the module\_type.
- 3. A given module instance in a given *art* process may make many data products of the same type. These are distinguished by giving each a unique instance name. An empty string is a valid instance name and in fact is the default. The other three fields must be non-empty strings.
- 4. The process name field holds the value of the process\_name parameter from the FHiCL file for the *art* job that created the data product.

The friendly name of a data type is a concept that *art* inherited from the CMS software suite. You will never need to write friendly names but you will need to recognize them. Knowing the following rules will be sufficient in most cases:

- 1. If a type is not a collection type, then its friendly name is the fully qualified name of the class. Fully qualified means that all namespace(s) and enclosing class names are included.
- 2. If a type is std::vector<T>, its friendly name is Ts; the mnemonic is that adding the letter "s" makes it plural.
- 3. If a type is std::vector<std::vector<T> >, its friendly name is Tss. And so on.
- 4. If a type is cet::map\_vector<T>, its friendly name is Tmv.

The full set of rules is given elsewhere.

Corollaries of the above discussion include:

• None of the four fields in a product name may contain an underscore character: otherwise the parsing of the name into its four fields is ambiguous.

If an art event-data file is populated by running several art jobs, each of which
adds some data products, then each art job in the sequence must have a unique
process\_name.

#### 15.4.3 Specifying a Data Product

To identify a data product, *art* requires that you specify the data type, module label and instance name fields (an empty string is a valid instance name). If the event contains exactly one data product that matches this specification, then *art* allows a wild card match on the process name field. If the event contains more than one data product that matches this specification, then *art* requires that you also specify the process name field. Note *art* allows a wild card match only on the process name field, not on the others.

The *art* Applications Programming Interface (API) treats the four parts of the data product name in two different ways. You use a template argument to tell *art* which data type you want. You use an object of type art::InputTag to specify the other three fields, module label, instance name and process name.

```
The header for art::InputTag is found in the file $ART_INC/art/Utilities/InputTag.h.
```

You can construct an input tag by passing it a string with the three fields separated by colons:

```
art::InputTag tag("ModuleLabel:InstanceName:ProcessName");
```

For this exercise, the full specification of the input tag is

```
art::InputTag tag("evtgen::exampleInput");
```

The double colon indicates that the instance name (which would come between the colons) is an empty string. The process name rarely needs to be specified, and in fact it is not needed in this exercise. It will be sufficient to specify the input tag as

```
art::InputTag tag("evtgen");
```

There are other constructors for art::InputTag and there are accessor methods that provide access to the individual fields. You can learn about these by looking at the header file but you will not use these features in this exercise.



#### 15.4.4 The Data Product used in this Exercise

The input files used for this exercise contain data products, one of which this exercise will use. This data product has the following attributes:

- o it has a data type of tex::GenParticleCollection
- o it is produced by a module with the label evtgen
- o its instance name is an empty string
- it is produced by an art job with the process name exampleInput.

# 15.5 Setting up to Run this Exercise

To run this exercise, you need to be logged in to the computer on which you ran Exercise 2 (in Chapter 10).

If you are continuing on from a previous exercise, you need to keep both your source and build windows open.

If you are logging back in, follow the instructions in Section 10.5 to reestablish your source and build windows.

In your source window, cd to your source directory. Then cd to the directory for this exercise and look at its contents:

```
ls
CMakeLists.txt
```

cd art-workbook/ReadGenParticles

```
readGens1.fcl ReadGens1\_module.cc
readGens2.fcl ReadGens2\_module.cc
readGens3.fcl ReadGens3\_module.cc
```

In this exercise you will run three modules that differ in only a few lines. The three source files use different syntax to accomplish the same thing. We recommend that, in most cases,

Listing 15.2: Output made using readGen1\_module.fcl

```
1 ReadGens1::analyze event: 1 GenParticles: 7
2 ReadGens1::analyze event: 2 GenParticles: 3
3 ReadGens1::analyze event: 3 GenParticles: 3
4 ReadGens1::analyze event: 4 GenParticles: 3
5 ReadGens1::analyze event: 5 GenParticles: 5
```

you use the syntax shown in the third version, ReadGens3\_module.cc; most of the exercises in the workbook will use this syntax. A description of the first two here serves as a pedagodical progression. You will likely see all three syntax versions in your experiment's code.

# 15.6 Running the Exercise

You will run the exercise from your build directory in your build window. The code is already built. To run this exercise, cd to your build directory and type the command: art -c fcl/ReadGenParticles/readGens1.fcl

This will make the usual *art* output, interspersed with the output made by readGens1.fcl. The output from this module is shown in Listing 15.2. For each event it prints the event number and the number of GenParticles in that event.

# 15.7 Understanding the First Version, ReadGens1

#### 15.7.1 The Source File ReadGens1 module.cc

Let's take a first look at the first module, shown in Figure 15.1. Read the explanations, then continue on to learn about *handles*.

```
#include ``toyExperiment/MCDataProducts/GenParticleCollection.h"
                                                       Notice the new
#include "art/Framework/Core/EDAnalyzer.h"
#include "art/Framework/Core/ModuleMacros.h"
                                                       include file at the top.
#include "art/Framework/Principal/Event.h"
#include <iostream>
#include <string>
namespace tex {
 class ReadGens1 : public art::EDAnalyzer {
  explicit ReadGens1(fhicl::ParameterSet const&);
  void analyze(art::Event const& event) override;
                              There is a new data member, gensTag_,
  art::InputTag gensTag_;
                              which is initialized in the constructor using a
 };
                              string value that is taken from the parameter set.
tex::ReadGens1::ReadGens1(fhicl::ParameterSet const& pset ):
 art::EDAnalyzer(pset),
 gensTag_(pset.get<std::string>("genParticlesInputTag")){
void tex::ReadGens1::analyze(art::Event const& event ){
 art::Handle<GenParticleCollection> gens; These two lines directly at left are the core of
 event.getByLabel(gensTag_,gens);
                                              the exercise: they set gens as a handle to the
                                              requested GenParticleCollection.
std::cout << "ReadGens1::analyze for event: " << event.id()
        << " the number of generated particles is: " << gens->size()
        << std::endl;
                         These std::cout lines print out the number of entries
                         in the data product: the same number as the number of
                         generated particles in the event.
DEFINE_ART_MODULE(tex::ReadGens1)
```

Figure 15.1: File listing for ReadGens1\_module.cc

Look at the line that begins art::Handle, which introduces the concept of a handle. As you work through the *art* workbook you will encounter several types of handles. All of the handle types behave like pointers with additional features:

- 1. They have safety features that make it impossible for your code to look at a pointee that is either not valid or not available.
- 2. They may also have an interface that lets you access metadata that describes the pointee.

The handle is an example of a broader idea sometimes called a *safe pointer*( $\gamma$ ) and sometimes called a *smart pointer*( $\gamma$ ).

The header for the class template art::Handle is found in the file \$ART\_INC/art/Framework/Principal/Handle.h. This file is automatically included by the include for Event.h.

The art::Handle line tells the compiler to default construct an object of type: art::Handle<GenParticleCollection>. The name of the default-constructed object is gens. A default-constructed handle does not point at anything and, if you try to use it as a pointer, it will throw an exception. A handle in this state is said to be invalid.

The following line calls <code>getByLabel</code>, which uses its first argument (<code>gensTag\_</code>) to learn three of the four elements of the name of the requested data product. It can deduce the fourth element, the data type, from the <code>type</code> of its second argument (<code>gens</code>): that is, it knows that it must look for a data product of type <code>tex::GenParticleCollection</code>. <code>art</code> has tools to compute the friendly name from the full class name, which is why you will never need to write a friendly name.

When this line is executed, the event object looks to see if it contains a data product that matches the request. There are three possible outcomes:

- 1. the event contains exactly one product that matches
- 2. the event contains no product that matches
- 3. the event contains more than one product that matches

In the first case, the event object will give the handle a pointer to the requested tex::GenParticleCollection; the handle gens can then be used as a pointer,

as is done in the second line of the std:cout section. When the handle has received the pointer, it is said to be in a valid state. In the second and third cases, the event object will leave the handle in its default-constructed state and, if you try to use it as a pointer, it will throw an exception.

If the event object finds exactly one match, it will also add two pieces of metadata to the handle. One is a pointer to an object of type art::Provenance, which contains information about the processing history of the data product. The second is an object of type art::ProductID; this is essentially a synonym for the four-field string form of the product name. Both of these will be illustrated in future exercises.

The third case bears one more comment: the developers of *art* made a careful decision that, except for the process name field, <code>getByLabel</code> will not have a notion of "best match". When you use <code>getByLabel</code> you must unambigously specify the data product you want or *art* will leave the handle in its default-constructed state.

If the getByLabel member function does not find the requested data product, e.g., if you run it on a different input file or if you misspell any of the fields in the input tag, the handle will be left in its default-constructed state. In this case, the gens->size() call will know that the handle is invalid and will throw an exception.

In all cases but one, *art*'s response to an exception is to attempt a graceful shutdown. The one unusual case is *ProductFotFound*, which is the exception thrown by an invalid handle when you try to use it as a pointer. In this case *art* will print an warning message, skip this module and attempt to run the remaining modules in the trigger paths and end paths.

It is possible to test the state of gens by using the member function gens.isValid(), which returns a bool. This not illustrated in the example because in most cases we recommend that you let *art* deal with this for you.



In the preceding discussion we did not mention that <code>getByLabel</code> is actually a *member function template*. There is no explicit template argument in the <code>event.getByLabel</code> line because the C++ template mechanism is able to deduce the template argument from the type of the second argument.

The art::Event object supports several other ways to request data products from the event, including a way to get handles to all data products that match a partial specification. This material is beyond the scope of this exercise.

Listing 15.3: A fragment from readGen1.fcl

#### 15.7.2 Adding a Link Library to CMakeLists.txt

ReadGens1\_module.so requires a link library that was not needed by previous exercises: \$TOYEXPERIMENT\_LIB/libtoyExperiment\_MCDataProducts.so. This library contains the object code for the classes and functions defined in the MCDataProducts subdirectory of the toyExperiment UPS product. In particular it contains object code needed by the data product tex::GenParticleCollection.

To add this library to the link list required a one line modification to CMakeLists.txt. If you compare this file to the corresponding file for the previous exercise, you will that CMakeLists.txt for this exercise contains one additional line:

```
$ { TOYEXPERIMENT_MCDATAPRODUCTS }
```

The string TOYEXPERIMENT\_MCDATAPRODUCTS is a cmake variable that was defined when you first ran the buildtool command. The translated value of this variable is the name of the required link library.

#### 15.7.3 The FHiCL File readGens1.fcl

There is only one fragment of readGens1.fcl that contains any new ideas. It is the fragment that configures the module label read, reproduced in Listing 15.3

In this fragment the parameter genParticlesInputTag specifies the input tag that identifies the data product to be read by this exercise.

We recommend that you always initialize input tags using parameters from the parameter set and that you never initialize them using strings defined within the code. This will allow you run the same module on data products with different input tags; this is a widely used



feature.



We further recommend that you do not provide a default value in the call to get the parameter value from the parameter set. This derives from a general recommendation that parameters affecting physics output should never have default values; the only parameters with default values should be those that control debugging and diagnostics.

## **15.8** The Second Version, ReadGens2

Version 2 of this exercise consists of the files ReadGen2\_module.cc and readGen2.fcl. To run this version, cd to your build directory and type the command:

```
art -c fcl/ReadGenParticles/readGens2.fcl
```

It will produce the same output as the previous two versions.

The only significant change from version 1 to version 2 is that lines

```
art::Handle<GenParticleCollection> gens;
event.getByLabel(gensTag_, gens);
```

have been replaced by the single (long) line:

```
art::ValidHandle<GenParticleCollection> gens =
    event.getValidHandle<GenParticleCollection>(gensTag_);
```

This version is a little verbose but that aspect will be addressed in version 3. Note that the class template art::Handle has been replaced by a new class template art::ValidHandle. Both class templates are defined in the same header file, \$ART\_INC/art/Framework/Principal/Handle.h.

The above line has functionality very similar to that of the two lines from version 1: the net result is that gens can be used as pointer to the requesed data product. It also has an interface to access the art::Provenance and the art::ProductID.

However, there are several signficant differences between art::Handle<T> and art::ValidHandle<T>:

Part

- Unlike an art::Handle<T>, which may be either valid or invalid, an art::ValidHandle<T> is guaranteed to be valid. It cannot be defaultconstructed.
- 2. A call to getValidHandle<T> will either return a properly constructed art::ValidHandle<T> or it will throw a ProductNotFound exception.
- 3. art::ValidHandle does not have an isValid() method.
- 4. Everytime that you use an art::Handle<T> as a pointer, it first checks that the pointer is valid. On the other hand, when you use an art::ValidHandle<T> as a pointer, no check is necessary; using an art::ValidHandle<T> as a pointer is as fast as using a bare pointer or a reference.

#### **15.9** The Third Version, ReadGens3

Version 3 of this exercise consists of the files ReadGen3\_module.cc and readGen3.fcl. To run this version, cd to your build directory and type the command:

#### art -c fcl/ReadGenParticles/readGens3.fcl

It will produce the same output as the previous two versions.

The only change from version 2 is that the call to getValidHandle has a slightly different syntax that provides the same behavior but is less verbose (shown here on two lines):

```
auto gens =
    event.getValidHandle<GenParticleCollection>(gensTag_);
```

This version uses a feature of C++ that is new in C++-11, the keyword auto. This keyword tells the C++ compiler to automatically determine the correct type for gens.

When you call the member function getValidHandle<T> the return type will always be art::ValidHandle<T>.

Version 3 is the version that we recommend you use but you can use any of the three. We introduced the version using the keyword auto as the last version because it is a handy

**Listing 15.4:** Warning message when the module label of a requested data product is misspelled.

```
1
2 %MSG
3 %MSG-w FailModule: ReadGens3:read 15-Jun-2014 10:00:24 CDT
    run: 1 subRun: 0 event: 5
4 Module failed due to an exception
5 ---- ProductNotFound BEGIN
6 getByLabel: Found zero products matching all criteria
7 Looking for type: std::vector<tex::GenParticle>
8 Looking for module label: genevent
9 Looking for productInstanceName:
10
11 ---- ProductNotFound END
```

shorthand when you know how to determine the correct type but it is very confusing if you do not know how to do so.

In future exercises we will use the pattern of version 3 regularly.

# 15.10 Suggested Exercises

```
Edit readGens3.fcl and supply the full input tag: genParticlesInputTag: "evtgen::exampleInput" Run art and observe that it works correctly.
```

Edit readGens3.fcl files and misspell the the requested module label, for example genParticlesInputTag: "genevent"

Run *art* and observe the warning messages, which should look like the message in Listing 15.4. Note that the warning message includes information about the module label (read), the module\_type (ReadGens3) and the fields of the requested data product. Observe that, for each event, *art* prints the warning message and continues with the next event.

Look at the last line of the *art* output and observe that *art* completed with status 0! This is because *art* treats ProductNotFound as a warning, not as an error that will initiate a shutdown.

You can reconfigure art so that a ProductNotFound exception will cause art to shutdown

**Listing 15.5:** Exception message for ProductNotFound when default Exceptions are disabled.

```
1
2 %MSG-s ArtException: PostCloseFile 15-Jun-2014 10:32:55 CDT PostEndRun
3 cet::exception caught in art
4 --- EventProcessorFailure BEGIN
5
    An exception occurred during current event processing
     ---- EventProcessorFailure BEGIN
6
7
       An exception occurred during current event processing
8
       --- ScheduleExecutionFailure BEGIN
9
         ProcessingStopped.
10
11
         ---- ProductNotFound BEGIN
12
           getByLabel: Found zero products matching all criteria
13
           Looking for type: std::vector<tex::GenParticle>
14
           Looking for module label: genevent
15
           Looking for productInstanceName:
16
17
           cet::exception going through module ReadGens3/read run: 1 subRun: 0 event: 1
18
         ---- ProductNotFound END
19
         Exception going through path end_path
20
       ---- ScheduleExecutionFailure END
21
     ---- EventProcessorFailure END
22
     cet::exception caught in EventProcessor and rethrown
23 ---- EventProcessorFailure END
24 %MSG
```

gracefully. To do this, edit your modified readGens3.fcl and add the following line inside the services parameter set:

```
scheduler : defaultExceptions : false
```

This line tells *art* that its response to all exceptions should be to attempt a gracefull shutdown. When you rerun *art* you should see output like that shown in Listing 15.5.

## **15.11** Review

In this chapter you have learned:

- 1. the type tex::GenParticleCollection
- 2. the four part identifier of data product and the class art::InputTag
- 3. the class templates art::Handle and art::ValidHandle

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- 4. how to get a handle to a data product, given its type and input tag
- 5. how to use a handle as a pointer to the requested data product
- 6. how to recognize a ProductNotFound warning message.
- 7. how to tell *art* to treat the ProductNotFound exception as a hard error that will initiate a graceful shutdown

## 16 Exercise 7: Making a Histogram

## 16.1 Introduction

One of the workhorse tools of HEP data analysis is ROOT. Among its many features are tools for data analysis, visualization, presentation and persistency. Indeed *art* uses ROOT for event-data persisency (i.e. to write event-data to files and read it back in). You have already had a strong hint of this because you have already seen the module\_types RootInput and RootOuput; and you have seen that the input files for these excercises all end in .root.

It is possible to use ROOT as a data processing framework but that will not be discussed here. Throughout the *art* workbook we use *art* as the data processing framework and we use ROOT as a toolkit.

ROOT supports a wide variety of classes that are used for data analysis, visualization and presentation. This exercise will use one of those classes, TH1D, as an example of how to use ROOT in the *art* environment. The class TH1D is used to create, fill and present 1-dimenstional histograms in which the content of each bin is represented by a double. You can use the other ROOT classes by simply following the pattern you learn for TH1D.

Much more information about ROOT is available from its own website, http://root.cern.ch/drupal.

In a typical *art* job, most of the modules run in that job, plus *art* itself, will all use ROOT. This presents an organizational problem: when you write a module, how can you be sure that your module's use of ROOT will not interfere with the use of ROOT by *art* or by other modules that are run in the same job? The solution to this problem is that users of *art* should interact with ROOT via an *art* service named the TFileService, which does

the necessary organizational work. In this chapter you will be introduced *art* services in general and the C++ TFileService in particular.

## 16.2 Prerequisites

Prerequisites for this chapter include all of the material in Part I (Introduction) and the material in Part II (Workbook) up to and including Chapter 15.

## 16.3 What You Will Learn

In this exercise you will learn,

- 1. What is the art::TFileService and what does it do for you.
- 2. How to configure the art::TFileService.
- 3. art::ServiceHandle
- 4. How to access ROOT via the art::TFileService
- 5. How to create and fill a ROOT TH1D histogram
- 6. How to use the interactive ROOT browser to view the histogram
- 7. How to run a CINT script to view the histogram and to write the histogram to a PDF file.
- 8. A naming convention used to distinguish event-data ROOT files from ROOT files that contains histograms, ntuples, trees and so on.

## 16.4 Setting up to Run this Exercise

To run this exercise, you need to be logged in to the computer on which you ran Exercise 2 (in Chapter 10).

If you are continuing on from a previous exercise, you need to keep both your source and build windows open.

If you are logging back in, follow the instructions in Section 10.5 to reestablish your source and build windows.

In your source window, cd to your source directory. Then cd to the directory for this exercise and look at its contents:

### cd art-workbook/FirstHistogram

ls

CMakeLists.txt FirstHist1\ module.cc

drawHist1.C firstHist1.fcl

The module FirstHist1\_module.cc is very much like the module ReadGens3\_module.cc from the previous exercise. The main difference is that it does not create any printout but it does create and fill a histogram of the number of generated particles in each event.

The file firstHist1.fcl is very much like the file readGens3.fcl from the previous exercise. The main difference is that it configures the TFileService.

The file drawHist1.C is a CINT script file that contains the commands to open a ROOT file, draw a histogram and write it to a PDF file.

The file CMakeLists.txt plays its usual role. Compared to the corresponding file for the previous exercise, it has two additional link libraries and a directive that drawHist1.C should not be built. You will see what this last item means in the full discussion of CMakeLists.txt.

## **16.5** The Source File FirstHist1\_module.cc

The C++ source code for this exercise is found in the file FirstHist1\_module.cc. Listing 16.1 shows a fragment of this file, showing the included headers and the declaration of the module class, FirstHist1. Compared to the file ReadGens3\_module.cc from the previous exercise, four new lines have been added to this file:

- 1. line 7, which includes the header for the art TFileService
- 2. line 9, which includes the header for the ROOT class TH1D
- 3. line 22, which declares the member function beginJob
- 4. line 29, which declares a new member datum, named hNGens\_, of type pointer to an object of type TH1D.

The name hNGens\_ was chosen because this pointer will eventually point at a histogram object that contains a histogram of the number of generated particles per event. The *art* workbook has adopted the style that all names for pointers to histograms begin with the lower case letter "h".

#### The two new headers can be found at:

```
$ART_INC/art/Framework/Services/Optional/TFileService.h
$ROOT_INC/TH1D.h
```

The conventions for including header files from ROOT differ from those for including header files from *art* and from **toyExperiment**. To remind you, *art* and the **toyExperiment** UPS product obey the following conventions:

- 1. the names of all classes and functions are inside a namespace, either art or tex.
- 2. the syntax to include a header file includes the package name as the first element of the path. This documents the package to which the header file belongs.

When ROOT was developed, namespaces were not supported robustly by many C++ compilers. Therefore the developers of ROOT chose a different set of conventions:

- 1. the names of all public root classes are in the global namespace (i.e. they are not part of a namespace that is defined by ROOT).
- 2. the names of all ROOT classes begin with a capital letter T followed by an upper case letter. This serves as a weak substitute for using a namespace.
- 3. When ROOT is distributed, a copy of every include file is put into one directory \$ROOT\_INC. The correct syntax to include a file from ROOT is to give the filename with any leading path elements. The clue that the file is a ROOT header file comes from the leading capital T.

**Listing 16.1:** The declaration of the class FirstHist1

```
1
2 #include "toyExperiment/MCDataProducts/GenParticleCollection.h"
3
4 #include "art/Framework/Core/EDAnalyzer.h"
5 #include "art/Framework/Core/ModuleMacros.h"
6 #include "art/Framework/Principal/Event.h"
7 #include "art/Framework/Services/Optional/TFileService.h"
8
9
  #include "TH1D.h"
10
11 #include <iostream>
12 #include <string>
13
14 namespace tex {
15
16
     class FirstHist1 : public art::EDAnalyzer {
17
18
     public:
19
20
       explicit FirstHist1(fhicl::ParameterSet const& );
21
22
       void beginJob() override;
23
       void analyze(art::Event const& event) override;
24
25
     private:
26
27
       art::InputTag gensTag_;
28
29
       TH1D* hNGens_;
30
31
     };
32
33
```

**Listing 16.2:** The implementation of the class FirstHist1

```
2
  tex::FirstHist1::FirstHist1(fhicl::ParameterSet const& pset ):
3
     art::EDAnalyzer(pset),
4
     gensTag_(pset.get<std::string>("genParticlesInputTag")),
5
     hNGens (nullptr) {
6
7
8
   void tex::FirstHist1::beginJob() {
9
10
     art::ServiceHandle<art::TFileService> tfs;
11
     hNGens = tfs->make<TH1D>( "hNGens",
12
               "Number of generated particles per event", 20, 0., 20.);
13
14 }
15
16 void tex::FirstHist1::analyze(art::Event const& event ) {
17
18
     auto gens = event.getValidHandle<GenParticleCollection>(gensTag_);
19
20
     hNGens_->Fill(gens->size());
21
22
```

Listing 16.2 shows the remaining section of the file FirstHist1\_module.cc. The new features in this file are,

- 1. line 5, which initializes hNGens\_ to have the value of a null pointer.
- 2. lines 8 through 14, the body of the beginJob member function, which creates an empty histogram.
- 3. line 20, which fills the histogram with the number of generated particles in the current event

The identifier nullptr, used in line 5, was added to the C++ core language in the 2011 Standard. It is the value of a pointer that points to nothing; in practice it has a value of zero. You will very likely encounter code written prior to the 2011 Standard. In such code you will see the equivalent of line 5 written in one of the following two ways: hNGens\_(0) or hNGens\_(NULL). In the second form, the value NULL is a C-Preprocessor MACRO variable that is defined to have a value of 0.

We strongly recommend that you use nullptr in your initializers. We strongly recommend that you never use the C-Preprocessor NULL.



## **16.5.1** Introducing art::ServiceHandle

Section 3.6.5 discussed the idea of *art* services; these are classes that provide some functionality (i.e. they provide a service) that can be used by any module or by other services. In this exercise, you will see your first example of an *art* service, the art::TFileService, which provides a bookkeeping layer to ensure that your use of ROOT does not interfere with others' use of ROOT.

Just as access to data products is provided by the class templates art::Handle and art::ValidHandle, access to services is provided by the class template art::ServiceHandle. Line 10 in Listing 16.2, tells the compiler to default construct an object, named tfs, of type art::ServiceHandle<art::TFileService>. The constructor of tfs will contact the internals of art and ask them to find a service of type art::TFileService. If art can find such a service, it will give the service handle a pointer to the service. If, on the other hand, art cannot find a service of the requested type, it will throw an exception and art will attempt a graceful shutdown.

Once a service handle has been constructed, the downstream code can use the service handle as a pointer to the pointee, in this case the art::TFileService.

The header file for art::ServiceHandle is found at: \$ART\_INC/art/Framework/Services/Registry/ServiceHandle.h It is automatically included by one of the files that are already included in FirstHist1 module.cc.

#### 16.5.2 Creating a Histogram

Lines 11 and 12 of Listing 16.2 use the  $\mathtt{art}: \mathtt{TFileService}$  to create a new histogram object of type <code>TH1D</code>. It returns a pointer to that object and the pointer is assigned to the member datum <code>hNGens</code> .

In the call to the member function template tfs->make, the type of object to be created is specified using a template argument. The non-template arguments are the arguments

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needed by a constructor of that type of object. You do not need to understand why things are done this way or how it all works. You just need to follow the pattern.

In the case of creating a TH1D, the five non-template arguments are:

- 1. The name by which ROOT will know this histogram; the art workbook has adopted the convention that this name will always be the name of the coresponding member datum, excluding the underscore.
- 2. The title that will be displayed when the histogram is drawn.
- 3. The number of bins in the histogram.
- 4. The lower edge of the lowest bin of the histogram.
- 5. The upper edge of the uppermost bin of the histogram.

ROOT defines that the low edge of a bin is within that bin, while the upper edge of a bin is part of the next bin up. Therefore the lower edge of the lowest bin is inside the histogram but the upper edge of the uppermost bin is outside of the histogram.

If you would like to learn more about the TH1D class you can look at its header file or you can read about it on the ROOT web site: http://root.cern.ch/root/html534/TH1D.html.

Where is the histogram created? The histogram is created in memory that is owned and managed by ROOT. ROOT also knows that when the job is finished, it should write the histogram to a ROOT output file so that you can inspect it at a later time. The name of the output file is specified in the FHiCL file for the *art* job; more on that later. In the following, we will call this file the *histogram output file* or simply the *histogram file*. Although *histogram files* often contain much more than just histograms, the name is in fairly common usage among the experiments that use *art*.

Just as file systems have the notion of directories and subdirectories (or folders and subfolders if you prefer), a ROOT file has the notion of directories and subdirectories that are internal to the ROOT file. If a module makes at least one histogram, then the TFileService will first create a new top level directory in the histogram file. The name of this top level directory is the name of the module label of the module that is creating the histogram. All ROOT objects that are created by that module will be created within this top level directory. When the contents of ROOT managed memory are written to the histogram file, this directory structure is preserved.

One of the rules of *art* is that, within one *art* job, each module label must be unique. This ensures that every module instance will have a uniquely named top level directory in the output ROOT file. It is this strategy that ensures that the histogram names of my module will never collide with the histogram names of your module.

The TFileService knows how to do its work in a way that does not interfere with *art*'s use of ROOT for input and output of event-data.

#### 16.5.3 Filling a Histogram

Line 20 of Listing 16.2 fills the histogram pointed to by hNGens\_ with the number of generated particles for this event.

If you look up the function prototype for TH1D::Fill you will see that it expects an argument that is a double. On the other hand, gens->size() returns an unsigned integer. One of the features of C++ is that it can automatically convert the unsigned integer to a double and pass that to the function.

#### 16.5.4 A Few Last Comments

All of the comments above about management of ROOT directories and writing histograms to files are also true for most other sorts of ROOT objects. In particular they are true for TTrees and TNtuples.

If you think carefully about FirstHist1\_module.cc you might wonder why there is no endJob member function containing a call to delete the histogram that was created in the beginJob member function. The answer is that when you create a histogram that is controlled by ROOT, then ROOT is responsible for calling delete at the right time.

If you talk to an HEP old-timer about creating histograms, they will probably call it "booking a histogram". This is language left over from a precursor to ROOT named HBOOK.

## 16.6 The Configuration File C++ firstHist1.fcl

The file firstHist1.fcl, shown in Listing 16.3, is very much like the file readGens3.fcl from the previous exercise. The most important new feature is at line 13, which configures the TFileService. This service has one required parameter, which is the name of the histogram file that contains the histograms, trees etc that are created by the *art* job.

If the required parameter is missing, or if the configuration for the TFileService is missing entirely, then the first attempt to get a service handle to the TFileService will throw an exception and *art* will attempt a graceful shutdown.

Another new feature is that this FHiCL file runs on the large input event-data file, inputFiles/input04\_data.root, which contains 1000 events.

#### 16.6.1 Two Kinds of ROOT files

By convention both the histogram files and the *art* event-data files end in .root. Even though but are ROOT files, the two types of files are structured very differently and are not in any way interchangable or interoperable. This can be confusing.

The *art* Workbook has adopted the convention that *art* event-data files always end in \_\_data.root. All other files ending in .root are histogram files.

Some experiments have adopted a similar convention while others have adopted precisely the opposite convention: files ending in \_hist.root are histogram files and all other files ending in .root are art event-data files.

## 16.7 The file CMakeLists.txt

The CMakeLists.txt file used for this exercise is shown in Listing 16.4. Compared to the corresponding file for the previous exercise, there are two new features.

- 1. Two link libraries have been added
- 2. There is a directive that no work should be done for files ending in . C

The two new link libraries are specified by

\${ART\_FRAMEWORK\_SERVICES\_OPTIONAL\_TFILESERVICE\_SERVICE}

**Listing 16.3:** The configuration file firstHist1.fcl

```
1
2 #include "fcl/minimalMessageService.fcl"
3
4 process_name : firstHist1
5
6 source : {
7
   module_type : RootInput
  fileNames : [ "inputFiles/input04_data.root" ]
9 }
10
11 services : {
message : @local::default_message
  TFileService : { fileName : "output/firstHist1.root" }
13
14 }
15
16 physics :{
17
   analyzers: {
18
      hist1 : {
19
                       : FirstHist1
       module_type
20
        genParticlesInputTag : "evtgen"
21
22
     }
23
24
    e1 : [ hist1 ]
25
     end_paths : [ e1 ]
26
27 }
```

Listing 16.4: CMakeLists.tex

```
file ( GLOB ROOT MACROS DO NOT BUILD
         RELATIVE ${CMAKE_CURRENT_SOURCE_DIR} *.C )
2
3
4 art_make(
5
    EXCLUDE ${ROOT MACROS DO NOT BUILD}
     MODULE LIBRARIES
6
7
       ${TOYEXPERIMENT_MCDATAPRODUCTS}
8
       ${ART_FRAMEWORK_CORE}
9
       ${ART_FRAMEWORK_PRINCIPAL}
10
       ${ART_PERSISTENCY_COMMON}
11
       ${ART FRAMEWORK SERVICES REGISTRY}
12
       ${ART_FRAMEWORK_SERVICES_OPTIONAL}
13
       ${ART_FRAMEWORK_SERVICES_OPTIONAL_TFILESERVICE_SERVICE}
14
       ${FHICLCPP}
15
       ${CETLIB}
16
       ${ROOT_HIST}
17)
```

and \${ROOT\_HIST}. These items are both cmake variables that were defined for you when you ran buildtool. The first variable translates to \$ART\_LIB/libart\_Framework\_Services\_Optional\_TFileService\_service.so and the second translates to a list of about 10 shared libraries from ROOT. This set of ROOT libraries is sufficient to let you use the most common ROOT classes, including histograms, ntuples and trees. If you want to use other ROOT classes you may need to further extend the link list. Consult a ROOT expert for details.

Many projects use the convention that files ending in .C contain code written in the C programming language. By default <code>cmake</code> will assume that files ending in .C follow this convention and, therefore, it will try to compile and link them. You will see in Section 16.11 that the file <code>drawHistl.C</code> is a script written in a language named CINT. Therefore <code>cmake</code> should simply ignore it.

Line 1 in CMakeLists.txt tells cmake to define a new cmake variable named ROOT\_MACROS\_DO\_NOT\_BUILD. This variable is the set of all filenames, from same directory as the CMakeLists.txt file, that end in .C. You do not need to understand how this line works.

Line 5 in CMakeLists.txt tells cmake that it should do nothing for all files that appear in the translation of the cmake variable ROOT\_MACROS\_DO\_NOT\_BUILD. Therefore cmake will ignore files ending in .C.

## 16.8 Running the Exercise

To run this exercise, cd to your build directory and type the command: art -c fcl/FirstHistogram/firstHist1.fcl

This module does not make any of its own printout. You should see the standard printout from *art*, including the final line saying that *art* will exit with status 0.

You should see that the *art* job created the file output/firstHist1.root. This is the file that earlier was called the histogram file.

## 16.9 Inspecting the Histogram File

In this section you will inspect the file output/firstHist1.root.

Before inspecting the this file, look again at fcl/FirstHistogram/firstHist1.fcl. Note that the module label of the FirstHist1 module is hist1.

To inspect the histogram you will remain in your build directory and you will run the interactive ROOT program, using the command root. This command was put into your path when you established your build environment. To perform this exercise:

#### 1. root -l

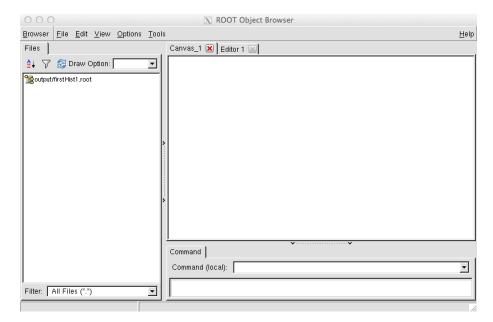
The command line option is a lower case letter "L". Some output and a new prompt will appear:

```
root [0]
Attaching file output/firstHist1.root as _file0...
root [1]
```

#### 2. At the root prompt, type the command

```
TBrowser* b = new TBrowser("Browser", _file0);
```

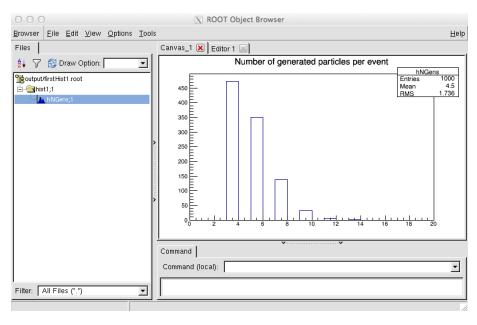
This will open a new window on your display; a screen capture of this window is shown in Figure 16.1. The text will refer to this window as the TBrowser wiindow.



**Figure 16.1:** Screen capture of the TBrowser window immediately after opening output/firstHist1.root.

- 3. In the left hand panel of the TBrowser window, you will see a ROOT file icon followed by the name of the file output/firstHist1.root. Click on this line.
- 4. This will create a new line in the left hand panel of the TBrowser window. The line contains a folder icon followed by the of a folder, hist1; 1. Click on this line.
- 5. This will create another new line in the left hand panel of the TBrowser window. This line contains a blue histogram icon and the name of a histogram hNGens; 1. Click on this line.
- 6. The histogram will appear in the right hand panel of the TBrowser window. Figure 16.2 shows a screen capture of the window with the histogram drawn.
- 7. To exit root, return focus to the build window and, at the root prompt, type .q ( a period followed by a lower case letter Q).
- 8. Another way to quit root is to click on the "Browser" pull-down menu on the top bar of the TBrowser window. From the menu select "Quit ROOT".

In step 4 you should have recognized the name of the folder, hist1; 1. Ignoring the



**Figure 16.2:** Screen capture of the TBrowser window after displaying the histogram hNGens; 1.

trailing; 1, it is the name of the module label used in firstHist1.fcl. In step 5 you should have recognized the name of the histogram, hNGens; ignoring; 1, it is the name that you gave the histogram when you created it.

About the ; 1 that ROOT has stiched onto hist1 and hNGens. ROOT calls these *cycle numbers*. They are part of a checkpointing mechanism that is beyond the scope of this exercise; if you ever see more than one cycle number for a ROOT object, the highest number is the one that you want. Consult the ROOT documentation for more details.

Now look at the histogram in the right hand panel of the TBrowser window. In the statistics box on the upper right you should see that it has 1000 entries, one for each event in the input file. You should also notice that only the odd bins are populated: this is because the generated events always contains three signal particles, plus a random number of pairs of background particles. The three signal particles are  $\varphi$  meson and the two kaons into which it decays. You should also recognize the title and the name that you set when you created the histogram. Finally you should recognize that the binning matches the binning you requested when you created the histogram.

## 16.10 A Short Cut: the browse command

The above description for viewing a ROOT file interactively requires a lot of tedious typing at step 2. The **toyExperiment** UPS product provides a command named browse that does the typing for you. To use this command:

#### browse output/firstHist1.root

Then follow the instructions from the previous section, starting at step 3.

When you created your *art* build in environment, the **toyExperiment** UPS product, put the command browse into your path. This command is implemented as bash script and you can find its definition using the following bash command:

#### type browse

browse is <...>/scripts/browse

where <...> will change from one site to the next; it will have the value of \$TOYEXPERIMENT\_DIR as defined at each site.

## 16.11 Using CINT Scripts

When you type a command at the ROOT prompt, you are typing commands in a ROOT-defined language called CINT. The name CINT is an acronym for C++ INTerpreter.\* As with many interpreted languages, you may write CINT commands in a file and execute that file as script.

It is a common convention that files that contain CINT scripts have a file type of . C. This convention is followed throughout the *art* workbook.

This exercise provides an example of a CINT script, drawHist1.C, which is shown in Listing 16.5. To use this script,

#### root -I drawHist1.C

This will open a window on your display, draw the histogram in that window and save the window to the PDF file output/NumberGenerated.pdf, which is shown as Figure 16.3. When the script is complete, it returns control to the root prompt in your

<sup>\*</sup> While CINT will correctly execute a lot of C++ code, there is legal C++ code that is not legal CINT and there is legal CINT that is not legal C++. There is also code that is legal in both C++ and CINT but which does subtley different things in the two environments. Therefore it is more correct to say that the CINT language shares a lot of syntax with C++, not that it is C++.

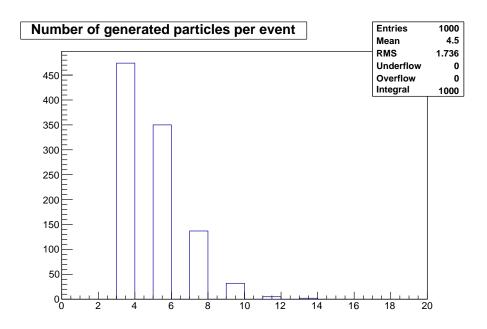


Figure 16.3: The figure made by running the CINT script drawHist1.C.

build window. At this prompt you can issue more C commands. To exit ROOT, type the command . q at the root prompt.

If you compare this figure to the histogram in Figure 16.2 you will see that there is difference in the statistics box in the upper right. In this figure, the name of the histogram is not shown but three new fields are shown: the number of entries below the lower limit (Underflow), the number of entries above the upper limit (Overflow) and the number of entries between the limits (Integral). The field named "Entries" is the sum of Integral plus Underflow plus Overflow. How to control what appears in the statistics box will be discussed in a few paragraphs.

It is beyond the scope of this writeup to describe all of the features in drawHist1.C so it will not discuss lines 6 and 9 of Listing 16.5. We will let comments in the code be a guide and give two additional hints. The object gROOT is a pointer to an instance of the class TROOT and the object gStyle is a pointer to an instance of the class TStyle. To find the documentation for these classes, see Section 16.11.1.

Line 14 tells ROOT what to draw in the statistics box in the upper right of every histogram. The comments in the code describe the mnemonics of the letter codes. The full set of letter

Listing 16.5: The CINT script drawHist1.C

```
1
2
3
4
     // With this you can reinvoke the script without exiting
5
     // root and restarting.
6
     gROOT->Reset();
7
8
     // Get rid of grey background (ugly for printing).
9
     gROOT->SetStyle("Plain");
10
11
     // Recommended content of statistics box:
12
     // Number of Entries, Mean, Rms, Underflows, Overflows,
13
     // Integral within limits
14
     gStyle->SetOptStat("emruoi");
15
16
     // Open the input file that contains histogram.
17
     TFile* file = new TFile( "output/firstHist1.root");
18
19
     // Get pointer to the histogram.
20
     TH1D* hNGens(0); file->GetObject("hist1/hNGens", hNGens);
21
22
     // Open a new canvas on the screen.
23
     TCanvas *canvas =
24
         new TCanvas("canvas", "Plots_from_Firsthist1.root");
25
26
     // "H9": draw outline histogram ("H") in high resolution mode (9)
27
     hNGens->Draw("H9");
28
29
     canvas->Update();
30
     canvas->Print("output/NumberGenerated.pdf");
31
32
```

codes is described on the ROOT web site: http://root.cern.ch/root/html534/TStyle.html On this page, navigate to the documentation for the member function SetOptStat.

Line 17 opens the input file. The ROOT type TFile is ROOT's interface to ROOT information that lives in a disk file; it allows a ROOT program to write ROOT objects to the file and read ROOT objects from the file. You can learn more about the class TFile from the ROOT web site http://root.cern.ch/root/html534/TFile.html.

Line 20 has two statements on it. The first declares hNGens as an object of type pointer to TH1D and initialzies it to 0 (ROOT does not support nullptr). The second asks the TFile to find a ROOT object named "hist1/hNGens", copy it from the file into memory and to set the pointer hNGens to point to that object. If ROOT cannot find the requested object, or if the type of the requested object does not match the type of the pointer, ROOT will leave hNGens with a value of 0. This is reminiscent of asking an art::Event to fill a an art::Handle.

Line 23 tells ROOT to open a new window on your display. The first argument is a name that ROOT uses internally to differentiate multiple canvases and the second argument is the title that will be drawn on the title bar of the window.

Line 27 tells ROOT to draw the histogram on the canvas. If, at line 20, ROOT was unable to properly set the pointer, then this line will produce an error message and return control to the root prompt in the build window.

Line 29 tells ROOT to flush its internal buffers and make sure that everything that is in the queue to be drawn on the canvas is actually drawn.

Line 30 tells ROOT to save the canvas by writing it to the file specified as the function argument. The format in which the file will be written is governed by the file type field in the filename, .pdf in this case. Many other formats are supported and a full list is available at: http://root.cern.ch/root/html534/TPad.html#TPad:Print

#### 16.11.1 Finding ROOT Documentation

The main ROOT web site is http://root.cern.ch/drupal. On the top navigation bar there is a title labeled "Documentation". Hover over this and a pull-down menu will appear. From this menu you can find links to Tutorials, How To's FAQs, a User Guide, a Reference

Manual and more. Some of the documentation is version dependent. The version of ROOT used by this version of the workbook is v5.34/18.

One possible starting point for learning ROOT is the ROOT Primer. You can find it by first going to the User's Guide page or you can follow the direct link: http://root.cern.ch/drupal/content/users-guide#primer

One of the most useful parts of the ROOT documentation suite is the Reference Guide, which can be reached from the pull down menu. The direct link to this page is: http://root.cern.ch/root/html534/ClassIndex.html

This section has a description of all of the members in each ROOT class.

## 16.12 Suggested Activities

### 16.12.1 Histogram Files are Overwritten

Suppose that you run *art* and make a histogram file. If you run *art* again, what happens? The answer is that it will overwrite the existing output file and replace it with the one created in the second job.

To illustrate this, rerun the exercise but tell *art* to only do 500 events:

art -c fcl/FirstHistogram/firstHist1.fcl -n 500

Inspect the output file and you will see that the histogram now has only 500 entries.

It is your responsibility not to overwrite files that you wish to keep.

### 16.12.2 Changing the Name of the Histogram File

You can change the name of the histogram file by editing the FHiCL file but you can also do so from the *art* command line by using the -TFileName option; the short form of this option is -T.

```
art -c fcl/FirstHistogram/firstHist1.fcl -TFileName output/anotherName.root -n 400 art -c fcl/FirstHistogram/firstHist1.fcl -T output/yetAnotherName.root -n 750
```

After each run, inspect the ouptut file and verify that the number of entries in the histogram matches the number of events requested on the command line.

## 16.12.3 Changing the Module Label

In firstHist1.fcl, change the name of the module label, hist1. Rerun the job and browse the histogram file. You should see that the name of the directory in the output file has changed to match the new module label.

### 16.12.4 Printing From the TBrowser

Use the browse command to open the histogram file and view the histogram. In the TBrowser window, click on the "File" button. This will open a pull-down menu; click on the line "Save As ...". This will open a dialog window that will let you save the histogram displayed on the canvas in a variety of formats, including .png, .gif, .jpg and .pdf.

## 16.13 Review

In this exercise you have learned

- 1. How to configure the TFileService
- 2. How to use an art::ServiceHandle to access the TFileService.
- 3. How to use the TFileService to create a histogram that will automatically be written to the output file.
- 4. Three different ways to view the contents of the histogram file: by launching a TBrowser by hand, by using browse command and by running a CINT script.
- 5. The convention used by the *art* Workbook to differentiate histogram files from *art* event data files.

# 17 Looping Over Collections

- 17.1 Prerequisites
- 17.2 What You Will Learn
- 17.3 Running the Exercise
- 17.4 Discussion
- 17.5 Suggested Activities

# 18 The Geometry Service

- 18.1 Prerequisites
- 18.2 What You Will Learn
- 18.3 Running the Exercise
- 18.4 Discussion
- 18.5 Suggested Activities

# 19 The Particle Data Table

- 19.1 Prerequisites
- 19.2 What You Will Learn
- 19.3 Running the Exercise
- 19.4 Discussion
- 19.5 Suggested Activities

# 20 GenParticle: Properties of Generated Particles

- 20.1 Prerequisites
- 20.2 What You Will Learn
- 20.3 Running the Exercise
- 20.4 Discussion
- 20.5 Suggested Activities

# Part III

# **Users Guide**

# 21 Obtaining Credentials to Access Fermilab Computing Resources

To request your Fermilab computing account(s) and permissions to log into the your experiment's nodes, fill out the form Request for Fermilab Visitor ID and Computer Accounts. Typically, experimenters that are not Fermilab employees are considered *visitors*. You will be required to read the Fermilab Policy on Computing.

After you submit the form, an email from the Fermilab Service Desk should arrive within a week (usually more quickly), saying that your Visitor ID (an identifying number), Kerberos Principal and Services Account have been created. You will need to change the password for both Kerberos and Services.

## 21.1 Kerberos Authentication

Your Kerberos Principal is effectively a username for accessing nodes that run Kerberos in what's called the FNAL.GOV *realm* (all non-PC Fermilab machines). \*

To change your Kerberos password, first choose one (minimum 10 characters with mixture of upper/lower case letters and numbers and/or symbols such as !, , #, \$, \$, &, \*, %). From your local machine, log into the machine using ssh or slogin and run the kpasswd command. Respond to the prompts, as follows:

\$ kpasswd <username>@FNAL.GOV

Password for username@FNAL.GOV: <--- type your current password here

<sup>\*</sup>The FERMI.WIN.FNAL.GOV realm is available for PCs.

```
New password: <--- type your new password here

New password (again): <--- type your new password here

Kerberos password changed.
```

Your Kerberos password will remain valid for 400 days.

## 21.2 Fermilab Services Account

The Services Account enables you to access a number of important applications at Fermilab with a single username/password (distinct from your Kerberos username/password). Applications available via the Services Account include SharePoint, Redmine, Service Desk, VPN and others.

To get your initial Services Account password, a user must first contact the Service Desk to get issued a first time default password. Once a default password is issued, users can access http://password-reset.fnal.gov/ to change it.

If you are not on-site or connected to the Fermi VPN, call the Service Desk at 630-840-2345. You will be given a one-time password and a link to change it.

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## 22 **git**

The source code for the exercises in the *art* workbook is stored in a source code management system called *git* and maintained in a repository managed by Fermilab. Think of *git* as an enhanced *svn* or (a VERY enhanced) *cvs* system. The repository is located at . You will be shown how to access it with *git*.

If you want some background on git, we suggest the Git Reference.

You will need to know how to install *git*, download the workbook exercise files initially to your system and how to download updates. You will not be checking in any code.

To install git on a Mac:

```
$ http://git-scm.com/download/mac
```

This will automatically download a disk image. Open the disk image and click on the .pkg file.

In your home directory, edit the file .bash\_profile and add the line:

```
$ export PATH=/usr/local/git/bin/:${PATH}
$ git clone ssh://p-art-workbook@cdcvs.fnal.gov/cvs/projects/art-workbook
```

and how to download updates as the developers make them:

```
$ git pull
```

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## 22.1 Aside: More Details about git

To bring your working copy of the workbook code up to date, you need to use git. Before describing the required git commands, we need to explain a little more about how git works and how the art-workbook team have chosen to use it. If you are familiar with git you can skip this section.

## 22.1.1 Central Repository, Local Repository and Working Directory

At any given time, there are three copies of the code that you need to be aware of, the central repository, your local clone of the central repository and the working copy of the code in your source directory.

- 1. The central git repository that contains all of the versions of the workbook is hosted by a machine named cdcvs.fnal.gov\* The art-workbook team updates this repository as it develops and maintains the exercises.
- 2. In section 10.4.1, in step 5b) you used the git clone command to make a copy of the central repository in your source directory. This clone contains a complete history of the development of art-workbook *as it existed at the time that you made the clone*. The local clone is found in the .git subdirectory of your source directory.
- 3. In section 10.4.1, in step 5d), you used the git checkout command to choose one of the tagged versions of art-workbook. This command looked into your local clone of the central repository, found all of the files in the requested version and put copies of them in the correct spot in the directory tree rooted at your source directory.



There are two other source code management systems that are widely used in HEP, cvs and svn. If you are familiar with either of these, git has an extra level: the concept of a local clone of the central respostory does not exist in those systems. That is, when are using cvs or svn and you want to switch to another version of the code, you need to contact the central repository but, when you are using git, you need only to contact your local clone of the central repository.

<sup>\*</sup> Originally this machine hosted only cvs repositories, hence its name. It now hosts cvs, svn and git repositories.

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To bring your working code up to date you need to do two steps:

- 1. Update your local clone of the central repository.
- 2. Checkout the new version from the local clone.

The discussion of the checkout has several cases. Each is discussed in one of the following sub-sections. It is possible for all four of these cases to occur on any given checkout.

## 22.1.1.1 Files that you have Added

When you worked on Exericse 2, you added some files to your working directories; for example you added the files Second\_module.cc and second.fcl. When you do the checkout of the new version, these files will remain in your working directory and will not be modifed; however the checkout command will generate some informational messages telling you that your working directories contain files that are not part of the checked out version.

You do not need to take any action; just be aware of the situation.

#### 22.1.1.2 Files that you have Modified

Another case occurs for files that have the following properties:

- 1. They were part of the old version.
- 2. You have modified them.
- 3. They have not been modified in the central repository since you cloned the repository.

For example, suppose that you modified first.fcl; it is very unlikely that this file would have been modified in the central repository after you cloned the repository.

In this case, the checkout command will issue a warning message to let you know that your working version contains changes that are not part of the release you checked out.

You do not need to take any action; just be aware of the situation.

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#### 22.1.1.3 Files with Resolvable Conflicts

Another case occurs for files that have the first two properties from the list in Section 22.1.1.2 but which have been modified in the central repository since you cloned the repository. This will happen from time to time when when we update exercises based on suggestions from users.

When this happens there are two cases, one of which is discussed here, while the other is discussed in the next sub-section.

If the two sets of changes (yours and those in the repository) are on different lines of the file git, will usually successfully merge these changes; git will then issue an warning message telling you want it has done.

It is your repsonsibility to identify these cases, understand the changes made in the repository and understand if git did the merge correctly.

#### 22.1.1.4 Files with Unresolvable Conflicts

The final case is a variant of the previous case; it occurs when git is unable to automatically merge conflicting changes. This will happen when the changes you made and the changes made in the repository affect the same line, or lines, of code. When git does not know how to merge the changes it will give up, add markup to the offending files to mark the conflict and issue an error message. This leaves the offending files in an unusable state and you must correct the conflicts, by hand, before continuing.

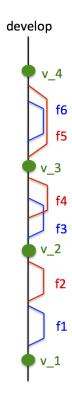
The art-workbook has been designed so that this should happen very, very rarely. Most readers should bookmark this spot for future reference and only read it when they need to.

#### **22.1.2** git Branches

git supports a concept known as *branches*. This is a very powerful feature that simplifies the task of having many developers collaboratively working on a single code base. Moreover, different experiments can choose to use branches in different ways; therefore a full description of branches is very open ended topic.

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Fortunately, to use the workbook you do not need to know very much about branches; all that you need to know is summarized in Figure 22.1, which shows a simplified view of the way that the art-workbook team uses git branches. In Figure 22.1 time starts at the bottom



**Figure 22.1:** A figure to illustrate the idea of git branches, as used in the Workbook; the figure is described in the text.

of the figure and runs upward. The art-workbook team has adopted the convention that the most up to date version of the art-workbook code will always be found by checking out a branch named *develop*. In Figure 22.1 the vertical line represents the develop branch.

At the earliest time represented in Figure 22.1, the develop branch existed in some state that the art-workbook team liked. So they tagged the develop branch with the name  $v_1$ , for version 1. Shortly afterwards, the art team needed to add some improvements. To do this they did:

1. Use the git pull command to make sure that their local copy of the respository is up to date.

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- 2. Check out the develop branch.
- 3. Start a new branch; in this example the new branch has the name £1, for "feature number 1".
- 4. Do all the development work on this branch. When they change files and commit their changes, the changes stay local to the branch.
- 5. Once the new code has been tested it is merged back into the develop branch.
- 6. Use the git pull command to make sure that their local copy of the respository is up to date; in this example, no one else has made.
- 7. Finally the developer must push their local copy of the respository to the central repository.

This is illustrated in Figure 22.1 by the blue line labelled £1. While the developer is working on £1, he can change back and forth between the develop branch and the £1 branch.

In Figure 22.1 the red line labelled £2 represents a second feature that is added to the code base following the same pattern as the first.

In this example, the development team decided to tag the develop branch after the £2 branch was merged back in; the tag was given the name v2; this is represented by the green filled circle in the figure.

The next items on the figure are the branches named £3 and £4. In this example, someone started with the develop branch and began work on the feature £3. A little later someone else (or maybe the same person) started with the develop branch and began work on the feature £4. The person starting work on £4 did so before the changes from £3 were merged back into the develop branch; therefore the two branches £3 and £4 both start from the same place, the v2 tag of develop. In this example, the next item on the timeline is that the developer of £3 commits their changes back to the develop branch. Sometime after that the developer of £4 merges their changes back. At this time the developer of £4 has the responsibility to check for conflicts that occurred during the merge and fix them; this may or may not require consultation with the author of £3.

After this, the develop branch is again tagged, this time with a version named v 3.

The next items on the timeline are the branches named £5 and £6. This example was

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included to show that it is legal for £6 both start and end during the time that £5 is alive.

Finally, the develop branch is tagged one more time, this time with the name  $v_4$ .

In Figure 22.1 consider a time when both branches £5 and £6 are active.

Fortunately, to use the workbook you do not need to know very much about branches. You really need to know only two things.

- 1. In the art-workbook repository, there is a branch named *develop*; this branch is the head of the project.
- 2. When the art-workbook team decides that a new stable version of the code is available, they checkout the develop branch and then start a new branch. This new branch, called a release branch has the same name as the version number of the release. For example, the code for version v0\_00\_13 is found in branch v0\_00\_13, and so on. New work, towards the next release, continues on the develop branch.

This is a bit of simplification but it captures the big ideas. Users of art-workbook should always work with one of the release branches; and they should always consult the documentation to learn which version of the code is matched to that version of the documentation.

Users of the art-workbook should never work in the develop branch; at any given time that branch may contain code that is still under development.

#### 22.1.3 Seeing which Files you have Modified or Added

At any time you can check to see which files you have modified and which you have added. To do this, cd to your source directory and issue the git status command. Suppose that you have checked out version  $v0\_00\_13$ , modified first.fcl and added second.fcl. The git status command will produce the following output:

```
$ git status

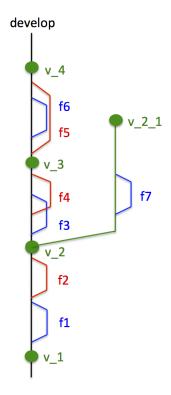
# On branch v0_00_13

# Changes not staged for commit:

# (use "git add <file>..." to update what will be committed)
```

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**Figure 22.2:** A figure to illustrate the idea of git branches, as used in the Workbook; the figure is described in the text.

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```
# (use "git checkout -- <file>..." to discard changes in
working directory)
#
# modified: first.fcl
#
# Untracked files:
# (use "git add <file>..." to include in what will be committed)
#
# second.fcl
no changes added to commit (use "git add" and/or "git commit -a")
```

You should not issue the git add or git commit commands that are suggested above.

In the rare case that you have neither modified nor added any files, the output of git status will be:

```
$ git status
# On branch v0_00_13
```

# 23 art Run-time and Development Environments

### 23.1 The art Run-time Environment

Your *art* run-time environment consists of:

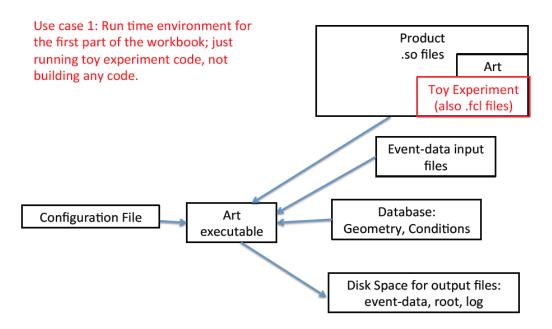
- o your current working directory
- all of the directories that you can see and that contain relevant files, including system directories, project directories, product directories, and so on
- o the files in the above directories
- the environment variables in your environment ( not sure how to say this nicely)
- o any aliases or shell functions that are defined

Figures 23.1, 23.2 and 23.3 show the elements of the run-time environment in various scenarios, and a general direction of information flow for job execution.

When you are running *art*, there are three environment variables that are particularly important:

- o PATH
- LD\_LIBRARY\_PATH
- FHICL\_FILE\_PATH

They are colon-separated lists of directory names. When you type a command at the command prompt, or in a shell script, the (bash) shell splits the line using whitespace and the

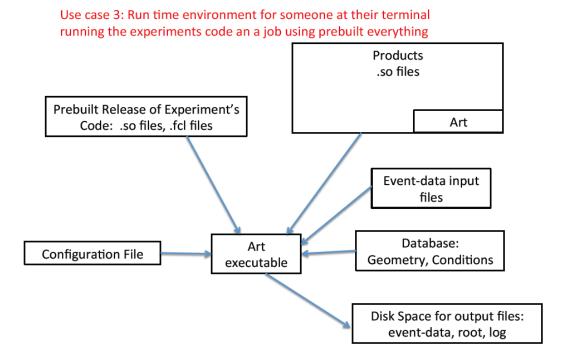


**Figure 23.1:** Elements of the *art* run-time environment, just for running the Toy Experiment code for the Workbook exercises

first element is taken as the name of a command. It looks in three places to figure out what you want it to do. In order of precedence:

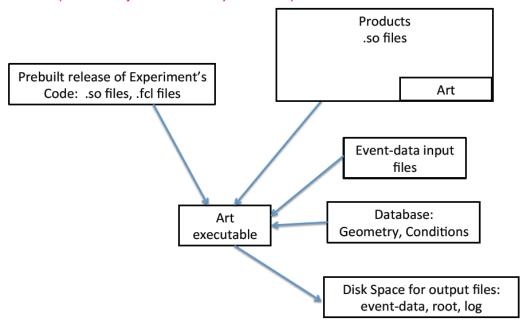
- 1. it first looks at any aliases that are defined
- 2. secondly, it looks for shell keywords in your environment with the command name you provide
- 3. thirdly, it looks for shell functions in your environment with that name
- 4. then it looks for shell built-ins in your environment with that name
- 5. finally, it looks in the first directory defined in PATH and looks for a file with that name; if it does not find a match, it continues with the next directory, and so on, followed by the paths defined in the other two variables.

Some parts of the run-time environment will be established at login time by your login scripts. This is highly site-dependent. We will describe what happens at Fermilab - consult your site experts to find out if anything is provided for you at your remote site.



**Figure 23.2:** Elements of the *art* run-time environment for running an experiment's code (everything pre-built)

Use case 4: Run time environment for someone running a production job with officially tracked inputs.



**Figure 23.3:** Elements of the *art* run-time environment for a production job with officially tracked inputs

When running the workbook, the interesting parts of your environment are established in two steps:

- o source a site-specific setup script
- o source a project-specific setup script

The Workbook, and the software suites for most IF experiments, are designed so that all site dependence is encoded in the site-specific setup script; that script adds information to your environment so that the project-specific scripts can be written to work properly on any site.

### 23.2 The art Development Environment

The development environment includes the run-time environment in Section 23.1 plus the following.

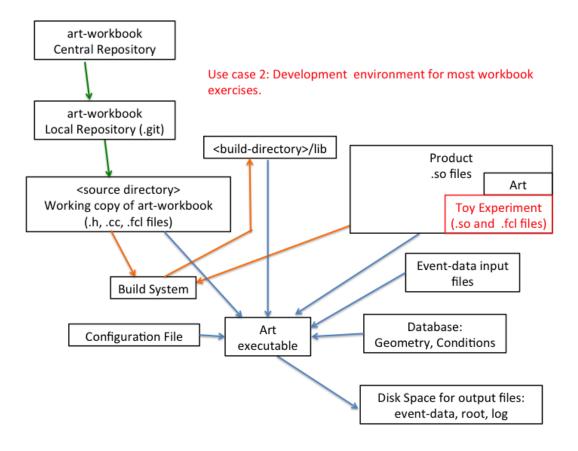
- the source code repository
- the build tools (these are the tools that know how to turn .h and .cc files in to .so files)
- additional environment variables and PATH elements that simplify the use of the above

Figures 23.4, 23.5 and 23.6 illustrate the development environment for various scenarios.

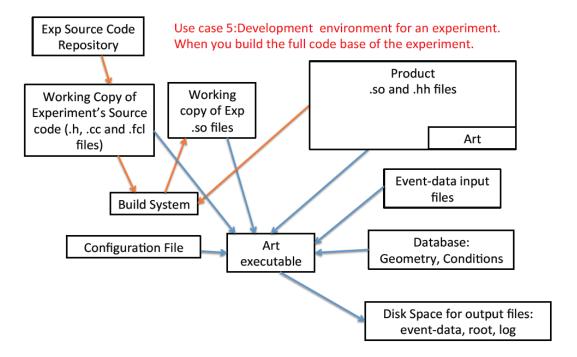
In some experiments the run-time and development environments are identical.

It turns out that there is no perfect solution for the job that build tools do. As a result, several different tools are widely used. Every tool has some pain associated with it. You never get to avoid pain entirely but you do get to pick where you will take your pain.

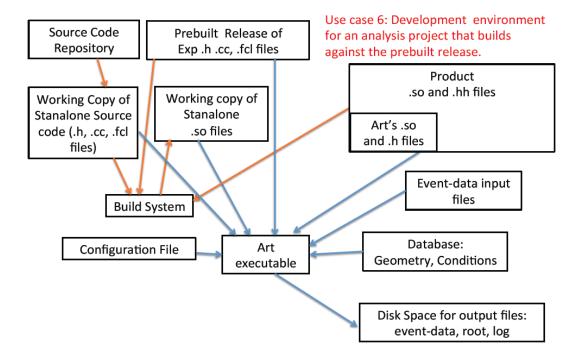
The workbook uses a build tool named **cetbuildtools**. Other projects have chosen make, cmake, scons and Software Release Tools (SRT). Here is something to watch out for: "build tools" written as two words refers generically to the above set of tools; but "build-tools" written as one word is the name of the executable that runs the build for **cetbuild-tools**.



**Figure 23.4:** Elements of the *art* development environment as used in most of the Workbook exercises



**Figure 23.5:** Elements of the *art* development environment for building the full code base of an experiment



**Figure 23.6:** Elements of the *art* development environment for an analysis project that builds against prebuilt release

## 24 art Framework Parameters

This chapter describes all the parameters currently understood by the *art* framework, including by framework-provided services and modules. The parameters are organized by category (module, service or miscellaneous), and preceded by a general introduction to the expected overall structure of an *art* FHiCL configuration document.

## 24.1 Parameter Types

The parameters are described in tables for each module. The type of a defined parameter may be:

```
TABLE: A nested parameter set, e.g., set: { par1: 3 }
SEQUENCE: A homogeneous sequence of items, e.g., list: [ 1, 1, 2, 3, 5, 8 ]
```

• STRING: A string (enclosing double quotes not required when the string matches [A-Za-z\_][A-Za-z0-9\_]\*). (Note: Identifiers when quoted do not function as special identifiers.) E.g.,

```
simpleString: g27
harderString: "a-1"
sneakystring1: "nil"
sneakystring2: "true"
sneakystring3: "false"
```

- COMPLEX: A complex number; e.g., cnum: (3, 5)
- NUMBER: A scalar (integer or floating point), e.g., num: 2.79E-8

```
o BOOL: A boolean, e.g.,
tbool: true
fbool: false
```

## 24.2 Structure of art Configuration Files

The expected structure of an art configuration file

Note, any parameter set is optional, although certain parameters or sets are expected to be in particular locations if defined.

```
# Prolog (as many as desired, but they must all be contiguous with only
# whitespace or comments inbetween.
BEGIN PROLOG
pset:
  nested_pset:
    v1: [ a, b, "c-d" ]
    b1: false
    c1: 29
  }
END_PROLOG
# Defaulted if missing: you should define it in most cases.
process_name: PNAME
# Descriptions of service and general configuration.
services:
{
  # Parameter sets for known, built-in services here.
  # ...
```

```
# Parameter sets for user-provided services here.
  user:
  {
  }
  # General configuration options here.
  scheduler:
  }
# Define what you actually want to do here.
physics:
  # Parameter sets for modules inheriting from EDProducer.
  producers:
    myProducer:
      module_type: MyProducer
      nested_pset: @local::pset.nested_pset
    }
  }
  # Parameter sets for modules inheriting from EDFilter.
  filters:
    myFilter: { module_type: SomeFilter }
  }
  # Parameter sets for modules inheriting from EDAnalyzer.
  analyzers:
  {
  }
```

```
# Define parameters which are lists of names of module sets for
  # inclusion in end_paths and trigger_paths.
  p1: [ myProdroducer, myFilter ]
  e1: [ myAnalyzer, myOutput ]
  # Compulsory for now: will be computed automatically in a future
  # version of ART.
 trigger_paths: [ p1 ]
  end_paths: [ e1 ]
}
# The primary source of data: expects one and only one input source
parameter set.
source:
# Parameter sets for output modules should go here.
outputs:
{
}
```

#### 24.3 Services

#### 24.3.1 System Services

These services are always loaded regardless of whether a configuration is specified.

#### 24.3.2 FloatingPointControl

These parameters control the behavior of floating point exceptions in different modules.

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 Table 24.1: art Floating Point Parameters

Enclosing	Parameter Name	Type	Default	Notes
Table Name				
services	floating_point_control	TABLE	{}	Top-level pa-
				rameter set for
floating point	setPrecisionDouble	BOOL	false	the service
floating_point_	setriccisionDouble	BOOL	Taise	
Control	reportSettings	BOOL	false	
	moduleNames	SEQUENCE	[]	Each module
				name listed
				should also
				have its own
				parameter
				set within float-
				ing_point_control.  One may also
				specify a mod-
				ule name of,
				"default" to
				provide default
				settings for the
due a duela	analdaDiaDa7anaE-	DOO!	folio a	following items:
<module-< td=""><td>enableDivByZeroEx</td><td>BOOL</td><td>false</td><td></td></module-<>	enableDivByZeroEx	BOOL	false	
name>	enableInvalidEx	BOOL	false	
	enableOverFlowEx	BOOL	false	
	enableUnderFlowEx	BOOL	false	

#### 24.3.3 Message Parameters

These parameters configure the behavior of the message logger (this is a pseudo-service – not accessible via ServiceHandle).

**Table 24.2:** art Message Parameters

Enclosing	Parameter Name	Type	Default	Notes
Table Name services	message	TABLE		Top-level pa-
				rameter set for the service
message				

#### 24.3.4 Optional Services

These services are only loaded if a configuration is specified (although it may be empty).

#### **24.3.5** Sources

#### **24.3.6** Modules

Output modules

## 25 Job Configuration in art: FHiCL

Run-time configuration for *art* is written in the Fermilab Hierarchical Configuration Language (FHiCL, pronounced "fickle"), a language that was developed at Fermilab to support run-time configuration for several projects, including *art*. For this reason, this chapter will need to discuss FHiCL both as a standalone language and as used by *art*.

By convention, the names of FHiCL files end in .fcl. Job execution is performed by running art on a FHiCL configuration file, which is specified via an argument for the -c option:

```
$ art -c run-time-configuration-file.fcl
```

See Figure ?? in Section 23.1 to see how the configuration file fits into the run-time environment.

The FHiCL concept of *sequence*, as listed in brackets [], maps onto the C++ concept of std::vector, which is a sequence container representing an array that can change in size. Similarly, the FHiCL idea of table, as listed in curly brackets {}, maps onto the idea of fhicl::ParameterSet. Note that ParameterSet is not part of *art*; it is part of a utility library *used* by *art*, FHICL-CPP, which is the C++ toolkit used to read FHiCL documents within *art*. FHiCL files provide the parameter sets to the C++ code, specified via module labels and paths, that is to be executed.

## 25.1 Basics of FHiCL Syntax

#### 25.1.1 Specifying Names and Values

A FHiCL file contains a collection of definitions of the form

#### name : value

where "name" is a parameter that is assigned the value "value." Many types of values are possible, from simple atomic values (a number, string, etc., with no internal whitespace) to highly structured table-like values; a value may also be a reference to a previously defined value. The white space on either side of the colon is optional. However, to include whitespace within a string, the string must be quoted (single or double quotes are equivalent in this case).

The fragment below will be used to illustrate some of the basics of FHiCL syntax:

```
# A comment.
// Also a comment.
name0 : 123
                         # A numeric value. Trailing comments
                         # work, too.
                          # Names can begin with underscores
name0 : 123
name00: "A quoted comment prefix, # or //, is just part of a
                                    # quoted string, not a comment"
name1:456.
                         # Another numeric value; whitespace is
                         # not important within a definition
name2 : -1.e-6
                         # A boolean value
name3 : true
NAME3 : false
                         # Other boolean value; names are case-
                         # sensitive.
name4 : red
                         # Simple strings need not be quoted
name5 : "a quoted string"
name6 : 'another quoted string'
name7 : 1 name8 : 2
                         # Two definitions on a line, separated by
                         # whitespace.
                         # Same as name9:3; newlines are just
name9
                         # whitespace, which is not important.
```

```
3
namea : [ abc, def, ghi, 123 ] # A sequence of atomic values.
                                # FHiCL allows heterogeneous
                                 # sequences, which are not,
                                 # however, usable via the C++ API.
                        # A table of definitions; tables may nest.
nameb :
   name0: 456
   name1: [7, 8, 9, 10]
   name2:
     name0: 789
 }
namec : [ name0:{ a:1 b:2 } name1:{ a:3 c:4 } ]
                        # A sequence of tables.
named : []
                       # An empty sequence
namee : {}
                        # An empty table
                       # An atomic value that is undefined.
namef : nil
abc : 1
                        # If a definition is repeated twice within
abc : 2
                       # the same scope, the second definition
def : [ 1, 2, 3 ]
                       # will win (e.g., "abc" will be 2 and
def: [4, 5, 6]
                       # "def" will be [4,5,6])
name : {
 abc : 1
 abc : 2
}
```

#### 25.1.2 FHiCL-reserved Characters and Identifiers

Several identifiers, characters and strings are *reserved to* FHiCL. What does this mean? Whenever FHiCL encounters a *reserved* string, FHiCL will interpret it according to the *reserved* meaning. Nothing prevents you from using these reserved strings in a name or value, but if you do, it is likely to confuse FHiCL. FHiCL may produce an error or warning, or it may silently do something different than what you intended. Bottom line: don't use reserved strings or symbols in the FHiCL environment for other than their intended uses.

The following characters, including the two-character sequence ::, are reserved to FHiCL:

```
, : :: @ [ ] { } ( )
```

The following strings have special meaning to FHiCL. They can be used as parameter values to pass to classes, e.g., to initialize a variable within a program, but their uses will not be fully described here because of subtleties and variations. As you work with C++ and FHiCL, the way to use them will become clearer.

true, false These values convert to a boolean

**nil** This value is associated with no data type. E.g. if a : nil, then a can't be converted

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to any data type, and it must be redefined before use

**infinity, +infinity, -infinity** These values initialize a variable to positive (the first two) or negative (the third) infinity

```
BEGIN_PROLOG, END_PROLOG ()
```

The first six strings (three lines) above function as identifiers reserved to *art* only when entered as lower case and unquoted; the last two strings (the last line) are reserved to *art* only when they are in upper case, unquoted and at the start of a line. Otherwise these are just strings. You may include any of the above reserved characters and identifiers in a "quoted" string to prevent them from being recognized as reserved to *art*.

#### 25.2 FHiCL Identifiers Reserved to art

FHiCL supports run-time configuration for several projects, not only for *art*. *art* reserves certain FHiCL names as identifiers that it uses in well-defined ways. (Other projects may use FHiCL names differently.) Within FHiCL files used by *art*, these FHiCL names obey scoping rules similar to C++. These identifiers appear in the FHiCL file with a scope, i.e.,

```
identifier : {
...
}
```

if they define a list of modules or a processing block, or with square brackets

```
identifier :[
...
]
```

if they define a list of paths.

The following is a list of the identifiers reserved to *art* and their meanings. In the outermost scope within a FHiCL file, the following can appear:

**process\_name** A user-given name to identify the configuration defined by the FHiCL file (it is recommended to make it similar to the FHiCL file name). This must appear at the top of the file. It may not contain the underscore character (\_).

**source** Identifies the data source, e.g., a file in ROOT format containing HEP events.

services Identifies ...

**physics** Identifies the block of code that configures the scientific work to be done on every event (as contrasted with the "bookkeeping" portions).

**outputs** List of output modules.

The following may appear within the physics scope:

producers Sets the list and order of producer modules; see Chapter 27

analyzers Sets the list and order of analyzer modules; see Chapter 28

filters Sets the list and order of filter modules; see Chapter 29

**trigger\_paths** List of producer and/or filter module paths; for each event, *art* executes all these module paths. The paths may only contain the module labels of producer and filter modules that are in the list of defined module labels. *art* can identify module labels that are common to several trigger\_paths and will execute them only once per event. The various paths within the trigger\_paths may be executed in any order.

**end\_paths** List of analyzer and/or output module paths; for each event, *art* executes all these module paths exactly once. The various paths within the end\_paths may be executed in any order.

The identifier process\_name is really only reserved to art within the outermost scope (but it would seem to be needlessly confusing to use process\_name as the name of a parameter within some other scope). The names trigger\_paths and end\_paths are artifacts of the first use of the CMS framework, to simulate the several hundred parallel paths within the CMS trigger; their meaning should be come clear after reading the remainder of this page.

## 25.3 Structure of a FHiCL Run-time Configuration File for art

Here is a sample FHiCL file called ex01.fcl that will do a physics analysis using the code in the *art* module  $Ex01\_module.so$  (the object file of the C++ source file

 $\texttt{Ex01}\_\texttt{module.cc}$ ). In this configuration, *art* will operate sequentially on the first three events contained in the source file inputFiles/input01\_data.root.

```
#include "fcl/minimalMessageService.fcl"
process_name : ex01
source : {
  module_type : RootInput
  fileNames : [ "inputFiles/input01_data.root" ]
  maxEvents : 3
}
services : {
 message : @local::default_message
}
physics :{
  analyzers: {
   hello : {
      module_type : Ex01
    }
  }
       : [ hello ]
  end_paths : [ e1 ]
}
```

Let's look at it step-by-step.

```
#include "fcl/minimalMessageService.fcl"
```

Similar to C++ syntax, this effectively replaces the '#include' line with the contents of the named file. This particular file sets up a messaging service.

```
process_name : ex01
```

The value of the parameter process\_name (ex01, here, the same as the FHiCL file name) identifies this *art* job. It is used as part of the identifier for data products produced in this job. For this reason, the value that you assign may not contain underscore (\_) characters. If the process\_name is absent, *art* substitutes a default value of "DUMMY."

```
source : {
  module_type : RootInput
  fileNames : [ "inputFiles/input01_data.root" ]
  maxEvents : 3
}
```

This source parameter describes where events come from. There may be at most one source module declared in an *art* configuration. At present there are two options for choosing a source module:

**module\_type : RootInput** art::Events will be read from an input file or from a list of input files; files are specified by giving their pathname within the file system.

**module\_type:** EmptyEvent Internally *art* will start the processing of each event by incrementing the event number and creating an empty art::Event. Subsequent modules then populate the art::Event. This is the normal procedure for generating simulated events.

Here RootInput is used; the data input file, in ROOT format, is assigned to the variable fileNames. The maxEvents parameter says: Look at only the first three events in this file. (A value of -1 here would mean "read them all.")

Note that if no source parameter set is present, *art* substitutes a default parameter set of:

```
source : {
  module_type : EmptyEvent
  maxEvents : 1
}
```

See the web page about configuring input and output modules for details about what other parameters may be supplied to these parameter sets.

```
services : {
  message : @local::default_message
}
```

Before starting processing, this puts the message logger in the recommended configuration.

```
physics :{
   analyzers: {
    hello : {
       module_type : Ex01
   }
}
```

In *art*, physics is the label for a portion of the run-time configuration of a job. It contains the "meat" of the configuration, i.e., the scientific processing instructions, in contrast to the more administrative or bookkeeping information. The physics block of code may contain up to five sections, each labeled with a reserved identifier (that together form a parameter set within the FHiCL language); the strings are *analyzers*, *producers*, *filters*, *trigger\_paths* and *end\_paths*. In our example it's set to analyzers.

The analyzers identifier takes values that are FHiCL tables of parameter sets (this is true also for filters and producers). Here it takes the value hello, which is defined as a table with one parameter, namely module\_type, set to the value Ex01. The setup defined a variable called LD\_LIBRARY\_PATH; art knows to match the value defined by the name module\_type to a C++ object file with the name Ex01\_module.so somewhere in the path defined by LD\_LIBRARY\_PATH.

We will expand on the physics portion of the FHiCL configuration in Section 25.5.

```
e1 : [ hello ] end_paths : [ e1 ]
```

## 25.4 Order of Elements in a FHiCL Run-time Configuration File for art

In FHiCL files there are very, very few places in which order is important. Here are the places where it matters:

- o A #include must come before lines that use names found inside the #include.
- A later definition of a name overrides an earlier definition of the same name.
- The definition of a name resolved using @local needs to be earlier in the file than the place(s) where it is used.
- Within a trigger path, the order of module labels is important.

Here is a list of *a few places* (of many) where order does not matter. This list is by no means exhaustive.

- Inside the physics scope, the order in which modules are defined does NOT matter for filters and analyzers blocks. These blocks define the run-time configurations of instances of modules.
- The five art-reserved words that appear in the outermost scope of a FHiCL file can be in any order. You could put outputs first and process\_name last, as far as FHiCL cares. It may be more difficult for humans to follow, however.
- Within the services block, the services may appear in any order.

Regarding trigger\_paths and end\_paths, the following is a conceptual description of how *art* processes the FHiCL file:

- art looks at the trigger\_paths sequence. It expands each trigger path in the sequence, removes duplicate entries and turns the result into an ordered list of module labels.
  The final list has to obey the order of each contributing trigger path, but there are no other ordering constraints.
- 2. It does the same for the end\_paths sequence but there is no constraint on order.
- 3. It makes one big sequence that contains everything in 1 followed by everything in 2.

- 4. It looks throughout the file to find parameter sets to match to each module label in the big list in 3.
- 5. It gives warning messages if there are left over parameter set definitions not matched to any module label in 3.
- 6. It then parses the rest of the physics block to make a "dictionary" that matches module labels to their configuration.

A conceptual description for the porcessing of services is as follows:

1. art first makes a list of all services, sorted alphabetically.



- 2. It makes a dictionary that matches service names to their parameter sets. A collorary is that service names must be unique within an *art* job.
- 3. *art* has some "magic" services that it knows about internally. It loads the .so file for each of them and constructs the services.
- 4. It loads the . so files for all of the services and calls their constructors, passing each service its proper parameter set.
- 5. It works through its list of modules in 5 it loads the .so and calls the constructor, passing the constructor the right parameter set.



6. It gives warning messages if there are left-over parameter set definitions not matched to any module label in 3.

When one service relies on another, things get a bit more complicated. If service A requires that service B be constructed first, then the constructor of service A must ask *art* for a handle to service B. When this happens, *art* will start to construct service A since it is alphabetically first. When the constructor of A asks for a handle to B, *art* will interupt the construction of service A, construct service B, and return to finish service A. Next, *art* will see that the next thing in the list is B, but it will notice that B has already been constructed and will skip to the next one.

Got that? Whew!

## 25.5 The *physics* Portion of the FHiCL Configuration

art looks for the experiment code in art modules. These must be referenced in the FHiCL file via module labels, which are just variable names that take particular values, as this section will describe. The structure of the FHiCL file – or a portion thereof – therefore defines the event loop for art to execute. The event loop, as defined in the FHiCL file, is collected into a scope labeled physics.

For a module label you may choose any name, as long as it is unique within a job, contains no underscore (\_) characters and is not one of the names reserved to *art*. In the sample physics scope code below, we define aProducer, bProducer, checkAll, selectMode0 and selectMode1 as module labels.

```
physics: {
producers : {
    aProducer: { module_type: MakeA }
    bProducer: { module_type: MakeB }
  }
 analyzers : {
    checkAll: { module_type: CheckAll }
  }
 filter : {
    selectMode0: {
      module_type: Filter1
      mode: 0
    selectMode1: {
      module_type: Filter1
      mode: 1
  }
```

The minimum configuration of a module is:

```
<moduleLabel> : { module_type : <ClassName> }
for example, in our code above:
aProducer: { module_type: MakeA }
```

aProducer is the module label and MakeA corresponds to a module of experiment code (i.e., an *art* module) named MakeA\_module.so, which in turn was built from MakeA\_module.cc. Since it falls within the scope producers, it must be a module of type EDProducer.

Let's take this a step farther, and assume that this EDProducer-type module MakeA accepts four arguments that we want to provide to *art*. The configuration may look like this:

```
moduleLabel : {
   module_type : MakeA
   pname0 : 1234.
   pname1 : [ abc, def]
   pname2 : {
       name0: {}
   }
}
```

This list under module\_type: MakeA represents parameters that will be formed into a fhicl::ParameterSet object and passed to the module MakeA as an argument in its constructor. pname0 is a double, pname1 is a sequence of two atomic character values, pname2 consists of a single table named name0 with undefined contents.

Note that *paths* are lists of module labels, while the two reserved names, trigger\_paths and end\_paths are lists of paths.

## 25.6 Choosing and Using Module Labels and Path Names

For a module label or a path name, you may choose any name so long as it is unique within a job, contains no underscore (\_) characters and is not one of the names reserved to *art* 

(see Section 25.2.

Any name that is a top-level name inside of the physics parameter set is either a reserved name or the name of a path.

It is important to recognize which identifiers are module labels and which are path names in a FHiCL file. It is also important to distinguish between a class that is a module and instances of that module class, each uniquely identified by a module label.

art has several rules that were recommended practices in the old framework but which were not strictly enforced by that framework. art enforces some of these rules and will, soon, enforce all of them:

- A path may go into either the trigger\_paths list or into the end\_paths list, but not both.
- A path that is in the trigger\_paths list may only contain the module labels of producer modules and filter modules.
- A path that is in the end\_paths list may only contain the module labels of analyzer modules and output modules.

Analyzer modules and the output modules may be separated into different paths; that might be convenient at some times but it is not necessary. On the other hand, keeping trigger paths separate has real meaning.

### 25.7 Scheduling Strategy in art

A set of scheduling rules is enforced in *art*. (Some of the details are remnants of compromises and conflicting interests with CMS.) One of the top-level rules in the scheduler is that all producers and filters must be run first, using the ordering rules specified below. After that, all analyzer and output modules will be run. Recall that analyzer modules and output modules may not modify the event, nor may they produce side effects that influence the behavior of other analyzer or output modules. Therefore, *art* is free to run analyzer and output modules in any order.

The full description of the scheduler strategy is given below:

- If a module name appears in the definition of a path name but it is not found among the list of defined module labels, FHiCL will issue an error.
- One each event, before executing any of the paths, execute the source module.
- On each event, execute all of the paths listed in the trigger\_paths.
  - Within one path, the order of modules listed in the path is followed strictly; at present there is one exception to this: see the discussion about the remaining issues
  - art can identify module labels that are in common to several trigger\_paths and will execute them only once per event. In the above example, aProducer and bProducer are executed only once per event.
  - The various paths within the trigger\_paths may be executed in any order, subject to the above constraints.
  - If a path contains a filter, and if the filter return false, then the remainder of the path is skipped.
  - The module name of a filter can be negated in path using, !moduleLabel; in this case the path will continue if the filter returns false and will be aborted if the filter returns true.
  - If the module label of a filter appears in two paths, negated in one path and not negated in the other, art will only run the instance of filter module once and will use the result in both places.
  - If a module in a trigger path throws, the default behaviour of art is to stop all processing and to shut down the job as gracefully as possible. Art can be configured, at run time, so that, for selected exceptions, it behaves differently. For example it can be configured to continue with the current trigger path, skip to the next trigger path, skip to the next event, and so on.
- o On each event, execute all of the paths listed in the end\_paths.
  - The module labels listed in end\_paths are executed exactly once per event, regardless of how many paths there are in the trigger\_paths and regardless of any filters that failed.

- If a module label appears multiple times among the end paths, it is executed only once. No warning message is given.
- Even if all trigger\_paths have filters that fail, all module labels in the end path will be run.
- End\_path is free to execute the modules in the end\_path in any order.
- If a module in the end\_path throws, the default response of art is to make a best effort to complete all other modules in the end path and then to shutdown the job in an orderly fashion. This behaviour can be changed at run-time by adding the appropriate parameter set to the top level .fcl file.
- One can ask that an output module be run only for events that pass a given trigger\_path; this is done using the SelectEvents parameter set,
- At present there is no syntax to ask that an analyzer module be run only for events that pass or fail some of the trigger paths. A planned improvement to art is to give analyzer modules a SelectEvents parameter that behaves as it does for output modules.
- If a path appears in neither the trigger\_paths nor the end\_paths, there is no warning given.
- If a module label appears in no path, a warning will be given.

In the above there is a lot of focus on which groups of modules are free to be run in an arbitrary order. This is laying the groundwork for module-parallel execution: *art* is capable of identifying which modules may be run in parallel and, on a multi-core machine, art could start separate threads for each module. At present both ROOT and G4 are not thread-safe so this is not of immediate interest. But there are efforts underway to make both of these thread-safe and we may one day care about module-parallel execution; our interest in this will depend a great deal on the future evolution of the relative costs of memory and CPU.

For simple cases, in which there is one trigger path with only a few modules in the path, and one end path with only a few modules in the path, the extra level of bookkeeping is just extra typing with no obvious benefit. The benefit comes when many work groups wish to run their modules on the same events during one art job; perhaps this is a job skimming off

many different calibration samples or perhaps it is a job selecting many different streams of interesting Monte Carlo events. In such a case, each work group needs only to define their own trigger path and their own end path, without regard for the requirements of other work groups; each work group also needs to ensure that their paths are added to the end\_paths and trigger\_paths variables. Art will then automatically, and correctly, schedule the work without redoing any work twice and without skipping work that must be done. This feature came for free with art and, while it imposes a small burden for novice users doing simple jobs, it provides an enormously powerful feature for advanced users. Therefore it was retained in art when some other features were removed.

### 25.8 Scheduled Reconstruction using Trigger Paths

Consider the following problem. You wish to run a job that has:

- Two producers MakeA\_module.cc and MakeB\_module.cc. You want to run both producers on all events.
- One analyzer module that you want to run on all events, CheckAll\_module.cc.
- You have a filter module, Filter1\_module.cc that has two modes, 0 and 1; the mode can be selected at run time via the parameter set.
- You wish to write all events that pass mode 0 of the filter to the file file0.root and you wish to write all events that pass mode 1 of the filter to file1.root

Here is code that would accomplish this:

```
process_name: filter1

source: {
    # Configure some source here.
}

physics: {
    producers : {
        aProducer: { module_type: MakeA }
```

```
bProducer: { module_type: MakeB }
 }
analyzers : {
   checkAll: { module_type: CheckAll }
 }
filter : {
   selectMode0: {
     module_type: Filter1
     mode: 0
   selectMode1: {
     module_type: Filter1
     mode: 1
  }
 mode0: [ aProducer, bProducer, selectMode0 ]
 model: [ aProducer, bProducer, selectModel ]
 analyzermods: [ checkAll ]
 outputFiles: [ out0, out1 ]
 trigger_paths : [ mode0, mode1 ]
 end_paths : [ analyzermods, outputFiles ]
}
outputs: {
 out0: {
  module_type: RootOutput
  fileName: "file0.root"
  SelectEvents: [ mode0 ] }
  }
 out1: {
```

```
module_type: RootOutput
fileName: "file1.root"
SelectEvents: { SelectEvents: [ mode1 ] }
}
```

Recall that the names process\_name, source, physics, producers, analyzers, filters, trigger\_paths, end\_paths and outputs are reserved to *art*. The names aProducer, bProducer, checkAll, selectMode0, selectMode1, out0 and out1 are module labels, and these are names of paths: mode0, mode1, outputFiles, analyzermods.

#### 25.9 Reconstruction On-Demand

#### 25.10 Bits and Pieces

What variables are known to *art*? physics (which has the five reserved identifiers: filters, analyzers, producers, trigger paths and end paths), what else? input file type RootInput

I know that trigger path are // different from end paths, they can contain different types of modules; // event gets frozen after trigger path.

art knows to match the value defi

ned by the name 'module\_name" to a C++ object fi

le with the name module\_name\_module.so" somewhere in the path defi

ned by LD LIBRARY PATH.

Further information on the FHiCL language and usage can be found at the mu2e FHiCL page.

# 26 Data Products

### 26.1 Overview

A *data product* is anything that you can add to an event or see in an event. Examples include the generated particles, the simulated particles produced by Geant4, the hits produced by Geant4, tracks found by the reconstruction algorithms, clusters found in the calorimeters and so on.

#### 26.2 The Full Name of a Data Product

Each data product within an event is uniquely identified by a four-part identifier that includes all namespace information. The four parts are separated by underscores:

```
DataType_ModuleLabel_InstanceName_ProcessName
```

*DataType* identifies the data type that is stored in the product. It is a "friendly" identifier in the way that its syntax has been standardized to deal with *collection* types, as follows:

- o If a product is of type T, then the friendly name is "T".
- If a product is of type mu2e::T, then the friendly name is "mu2e::T".
- If a product is of type std::vector(mu2e::T), then the friendly name is "mu2e::Ts".
- $\circ$  If a product is of type std::vector $\langle$ std::vector $\langle$  mu2e::T $\rangle$  $\rangle$ , then the friendly name is "mu2e::Tss".
- If a product is of type cet::map\_vector(mu2e::T), then the friendly name is "mu2e::Tmv".

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See below for a discussion about where underscores may not be used; this example is safe because of the substitution of "mv" for map\_vector.

*ModuleLabel* identifies the module that created the product; this is the module label, which distinguishes multiple instances of the same module within a produces. It is *not* the class name of the module.

*InstanceName* is a label for the data product that distinguishes two or more data products of the same type that were produced by the same module, in the same process. If a data product is already unique within this scope, it is legal to leave this field blank. The instance label is the optional argument of the call to "produces" in the constructor of the module (xxxx below):

```
produces<T>("xxxx");
```

*ProcessName* is the name of the process that created this product. It is specified in the FHiCL file that specifies the run-time configuration for the job (shown as *ReadBack02* below):

```
process_name : ReadBack02
```

Because the full name of the product uses the underscore character to delimit fields, it is forbidden to use underscores in any of the names of the fields. Therefore, none of the following items may contain underscores:

- o the class name of a class that is a data product; the exception is the cet::map\_vector template; when creating the friendly name, *art* internally recognizes this case and protects against it
- the namespace in which a data product class lives
- o module labels
- data product instance names
- o process names

It is important to know which names need to match each other; see Section 32.1.

# **27 Producer Modules**

# 28 Analyzer Modules

Analyzer modules request data products, do not create new ones; make histograms, etc.

•

An analyzer interface looks like the following.

```
class EDAnalyzer {
    // explicit EDAnalyzer(ParameterSet const&)

    virtual void analyze(Event const&) = 0
    virtual void reconfigure(ParameterSet const&)

    virtual void beginJob()
    virtual void endJob()
    virtual bool beginRun(Run const &)
    virtual bool endRun(Run const &)
    virtual bool endRun(SubRun const &)
    virtual bool endSubRun(SubRun const &)

    virtual void respondToOpenInputFile(FileBlock const& fb)
    virtual void respondToCloseInputFile(FileBlock const& fb)
    virtual void respondToOpenOutputFiles(FileBlock const& fb)
    virtual void respondToOpenOutputFiles(FileBlock const& fb)
    virtual void respondToCloseOutputFiles(FileBlock const& fb)
}
```

# 29 Filter Modules

Filter modules request data products and can alter further processing using return values

A filter interface looks like the following.

```
class EDFilter {
    // explicit EDFilter(ParameterSet const&)

    virtual bool filter(Event&) = 0
    virtual void reconfigure(ParameterSet const&)

    virtual void beginJob()
    virtual void endJob()
    virtual bool beginRun(Run &)
    virtual bool endRun(Run &)
    virtual bool endSubRun(SubRun &)
    virtual bool endSubRun(SubRun &)

    virtual void respondToOpenInputFile(FileBlock const& fb)
    virtual void respondToCloseInputFile(FileBlock const& fb)
    virtual void respondToOpenOutputFiles(FileBlock const& fb)
    virtual void respondToCloseOutputFiles(FileBlock const& fb)
    virtual void respondToCloseOutputFiles(FileBlock const& fb)
}
```

# 30 art Services

This chapter describes what services are, what types of services are recognized, how they are used, how they are constructed.

#### 30.1 About Services

i

art Services are introduced in Section ?? . A service in the context of art is a class for which an instance is configured at runtime by a FHiCL configuration and made available for use by modules and other services, and whose lifetime and configuration are managed by art. A service provides modules access to resources or information that is outside the purview of the modules, and that is valid for some aggregation of events, subRuns or runs, or over some time interval. Services may also be used to provide certain types of utility functions.

How services fit into *art* during job processing. *art* first pulls in the parameter set from the user's FHiCL file, instantiates the services called for, calls the constructor for each module, then begins the event loop.

Services can be provided by three different sources. Some services are provided by *art* to implement internal functions used by both *art* itself and experiment code, e.g., the message service. Other services implement functions that are not internal to *art* but are used by most experiments, for example TFileService, which manages a secondary ROOT-format file for histograms, and and, which provide interfaces to geometry and calibration information, respectively. Users (experiments and/or individual experimenters) can add services to *art*, too.

Services can be publicly available, i.e., callable by user-defined classes, or callable by other services or by *art* itself. Services that are publicly available typically have a header (.h) file, whereas those intended to be called by *art* itself do not. A service may *register* functions with *art*; a registered function can then be called at appropriate moments directly by the *art* framework, e.g., before or after beginRun, endSubRun and so forth.

Services can be made available for environments that support parallelism. More in Section 30.6.1.

A service should *not* provide "backdoor" communication between modules of physics data or to/from anything that ought to be stored in an event, subRun or run, and/or anything that requires provenance tracking, since the provenance would not be captured.

To access a given service, the calling entity requests a *service handle* for it; this is a type of smart pointer that ... Depending on what a service does and what external system(s), if any, it needs to access, a *service interface* may be required. .

Several types of standard *art* services exist:

- TFile: Controls the ROOT directories (one per module) and manages the histogram file.
- o Timing: Tracks CPU and wall clock time for each module for each event
- Memory: Tracks increases in overall program memory on each module invocation
- FloatingPointControl: Allows configuration of FPU hardware "exception" processing
- RandomNumberService: Manages the state of a random number stream for each interested module
- srvc MessageFacility: Routes user-emitted messages from modules based on type and severity to destinations

## 30.2 Service Handles

To use a service, your code must point to it. Your FHiCL file must point to the service(s) you need and include its configuration, e.g., for TFileService:

art Documentation

```
1 #include "fcl/minimalMessageService.fcl"
2 process_name : ex06
3 ...
4 services : {
5     ...
6    TFileService : { fileName : "output/ex06.root" }
7  }
8    ...
```

The .h or .cc file's declaration of your module (class) under "private" must declare a service handle to the service. This shows part of the code in Ex06\_module.cc in which a service handle to the TFileService is declared with the name tfs\_. The is in the namespace art.:

```
class Ex06 : public art::EDAnalyzer {
public:
    explicit Ex06(fhicl::ParameterSet const&);

...

private:
    art::ServiceHandle<art::TFileService> tfs_;
```

The class's constructor calls/declares? it as a member function of the class, with no arguments:

```
1 tex::Ex06::Ex06(fhicl::ParameterSet const& pset ):
2   gensTag_(pset.get<std::string>("genParticlesInputTag")),
3   tfs_(),
4   hNGen_(nullptr) {
5 }
```

# 30.3 Implementing Simple Services

A class that is intended as a service must be declared as a service to the *art* framework and its implementation defined. The way to do this depends on whether the service is provided by *art* or not, and whether it requires an interface.

An art-provided service such as Timing or SimpleMemoryCheck should be configured for use in this way. In the FHiCL file

```
services.SimpleMemoryCheck: { }
```

Part

Services not provided by art must be configured to be nested under services.user:

```
services.user.MyService: { }
```

Services implementing an interface require yet another syntax for configuration. In services or services.user there must be the line (with the appropriate ServiceName):

```
InterfaceName: { service_provider: ServiceName }
For example:
services.user.CatalogServiceInterface: {
    service_provider: IFCatalogInterface
}
```

# 30.4 Configuring a Service

First, a service needs to be configured for use. In a FHiCL file, this would look like (shown for the service TFileService):

## 30.5 Accessing a Service

Services are accessed via an instantiation of the art::ServiceHandle template. E.g., to access the TFileService from your module, use this syntax:

```
art::ServiceHandle<art::TFileService> h;
```

Treat the handle so created as any other smart pointer.

If you intend to access one service from a second service, you should obtain a ServiceHandle for the second service from the constructor of the first, even if you don't plan to use it there. This ensures the correct order of construction (and destruction) of the services, which otherwise would be alphabetical by class name.

Any library containing code that uses a service should link explicitly to that service.

Note that obtaining an art::ServiceHandle is relatively expensive in terms of performance, and should not be done in tight loops. It is recommended to cache the handle in your module class scope, at the class or module entry point (i.e., analyze(), filter()).



To avoid problems when parallelism is in operation, do not cache a bare pointer to a service at a greater-than-module-entry-point scope.

De-referencing the handle, on the other hand, while not zero-cost, is relatively cheap. As for certain types of services, *art* runs a check on the handle to verify that the service referenced is appropriate for the context in which the handle will be used. If necessary, the handle is "re-seated" for the new context.

# 30.6 Writing a Service

ServiceMacros.h

The code in the .cc file implementing a service must begin with a call to the *art service macro*: What does ServiceMacros do?

#include "art/Framework/Services/Registry/ServiceMacros.h"

Part

If the service implements an interface, the interface definition takes the form of a header file containing a class definition (declaration?) and interface declaration. Clarify 'define' vs 'declare' vs 'implement'; I don't think we're using them consistently (sec 6.6.4 says in ptest.cc a line declares the variable p0 even though ptest.h exists)

An access interface looks like the following (shown for the service TFileService):

An interface implementation must inherit publicly from the interface, and declare and define itself as an implementation of that interface.

```
#include "art/Framework/Services/??/???.h"
```

#### **30.6.1** Declaring and Defining Services

In general, DECLARE macro invocations should be in the header file, if it exists. DEFINE macro invocations should be in the implementation.

A service may have one of three scope indicators:

- **LEGACY** All currently defined services are "legacy" in scope, including *art*-defined services. You should (for now) define all your services to be "LEGACY". The presence of even one LEGACY service in the configuration precludes the simultaneous processing of multiple events ("multi-schedule" operation).
- **GLOBAL** Defining a service as "global" in scope gives a guarantee of thread saftey: the service may be called simultaneously from anywhere at any time and continue to behave correctly.
- **PER\_SCHEDULE** These services may be instantiated multiple times if the *art* framework is configured for multi-schedule operation.

For example, this *declaration* would be included in ServiceMacros.h:

```
DECLARE_ART_SERVICE (myexpt::MyService, LEGACY)
```

whereas this *definition* would be in the implementation file (after the include statement for the header):

```
DEFINE_ART_SERVICE (myexpt::MyService)
```

Next you declare an interface; follow that by declaring then defining a service that will implement the interface:

Next, include the constructor:

```
MyService(fhicl::ParameterSet const &, art::ActivityRegistry &);
```

Via, the service is provided at construction time with access to its configuration parameters and to the <code>source:art/Framework/Services/Registry/ActivityRegistry.h</code> for the purposes of registering for callbacks at any of a large number of synchronization points in the execution of the *art* framework.

To register for callbacks, invoke a watch() function on the signal data member of ActivityRegistry that corresponds to the point in the execution of the *art* framework at which your function is to be called. For instance:

```
aReg.sPreBeginRun.watch(&MyService::preBeginRun, *this);
```

The function you provide must have the signature specified for that signal. The first template parameter of the signal specifies the calling order of registered callbacks; the second is the return type of the callback function; any subsequent parameters specify the ordered parameters of the callback function. The provided function may be any of:

- o a free function, bound or unbound
- o a const or non-const member function of your service or of any other class, bound or (if you provide an object with which to bind it) unbound
- o a function whose operator () signature matches the required signature

A service plugin is compiled into a plugin recognizable to the *art* framework by making a library whose name is lib[dir1\_dir2\_]ClassName\_service.so.

#### 30.7 Service Interfaces

<Interface> is an interface that can point to service MyService and YourService and TheirService. 'something' is a name that a user module can use to access any of these services. Is it an implementation? How does that work? What's the syntax for that?

A service interface defines a base class that is registered to *art* as such, thereby allowing a service to inherit from this base class/interface and declare itself to the *art* framework. Subsequently, *art* can be configured at runtime via FHiCL to use a given service that implements a particular interface. External users of that interface (i.e., other services, or on occasion the *art* framework itself) access the given service via the interface without ever knowing (or needing to know) exactly which (other) service implements it.

Art can be configured and/or a user module can be configured to use an interface?

A service that requires an interface must have a header? Or only if used by user module? And the header must include (what in the parens is mandatory?):

DECLARE\_ART\_SERVICE\_INTERFACE(myexpt::Interface,LEGACY)

# 31 art Input and Output

# 31.1 Input Modules

#### 31.1.1 Configuring Input Modules to Read from Files

When reading from an existing file, *art* allows you to select the input files, the starting event, the number of events to read, etc., either from the command line or from the FHiCL file. If a particular quantity is controlled from both the command line and the FHiCL file, the value on the command line takes precedence.

The following code fragment tells *art* to read event data from the file of type "ROOT," named "file01.root" and to start at the beginning of the file. A value of "-1" for maxEvents tells *art* to read events until the end of file is reached:

To tell *art* to read 100 events, or until the end of file, which ever comes first, change the parameter maxEvents to 100. This also shows how to specify a list of input files:

The number of files in the list of input files is arbitrary. The following fragment tells *art* to skip the first two events (and thus start with the third):

The fragment below shows some other parameters that can be included in the source pa-

**Listing 31.1:** Reading in a ROOT data file

```
1 source :{
2   module_type : RootInput
3   fileNames : [ "file01.root" ]
4   maxEvents : -1
5 }
```

**Listing 31.2:** Reading in a ROOT data file

```
1 source : {
2  module_type : RootInput
3  fileNames : [ "file01.root", "file02.root", "file03.root" ]
4  maxEvents : 100
5 }
```

**Listing 31.3:** Reading in a ROOT data file

```
1 source : {
2   module_type : RootInput
3   fileNames : [ "file01.root", "file02.root", "file03.root" ]
4   maxEvents : 100
5   skipEvents : 2
6 }
```

#### rameter set:

The parameters whose names start with *first* specify that the first event to be processed will be the first event that has an EventID greater than or equal to the specified event. If one of the first\* parameters is not specified, it takes a default value of -1 and is excluded from the comparison.

If a file of unsorted events is read in, *art* will, by default, present the events for processing in order of increasing event number. As a corollary to this, the output file will contain the events in sorted order. This sorting occurs one input file at a time; *art* does not sort across file boundaries in a list of input files. If the noEventSort parameter is set to *true*, the sorting is disabled, which will, in most cases, yield a minor performance improvement.

I have not yet learned the precise meaning of the skipBadFiles and the fileMatchMode

Listing 31.4: Reading in a ROOT data file

```
1
     firstRun
                          : 0
2
                          : 0
     firstSubRun
3
     firstEvent
                          : 0
4
     noEventSort
                          : false
5
                          : false
     skipBadFiles
6
     fileMatchMode
                          : "permissive"
7
     inputCommands
```

**Listing 31.5:** Reading in a ROOT data file

```
1 source :{
2  module_type : EmptyEvent
3  maxEvents : 200
4 }
```

#### parameters.

The inputCommands parameter tells *art* to delete certain data products from the copy of the event in memory after reading the input file. In other words, the input file itself is not modified but data products are removed from the copy of the event in memory before any modules are called. The syntax of this language is the same as for outputCommands, described.

In the pre-art versions of the framework, there were methods to select ranges of events or ranges of SubRuns. This is not yet working in art; the art developers will add this feature back once we decided exactly what we mean by "ranges of events".

Specifying Many Input Files In the pre-art, python based, configuration language, the standard syntax to initialize a list of input files was limited to 255 files, after which an alternate syntax was required. This is no longer necessary; the length of a fhicl list is limited only by available memory.

#### **Empty Source**

In many simulation applications one wishes to start with an empty event, run one or more event generators, pass the generated particles through the Geant4, and so on. In art the first step in this chain is accomplished using a source module named EmptySource, as follows:

Instead of reading event-data from a file, the empty source increments the event number and presents an empty event to the modules that will do the work. One may configure EmptySource to specify the EventId of the first event, to specify the maximum number of events in a SubRun or SubRuns in a run.

The last option tells art to reset event numbers to start at 1 whenever art starts a new SubRun begins; this is the default behaviour and is opposite to the behaviour we inherited from CMS.

**Listing 31.6:** Reading in a ROOT data file

```
1 source : {
2
   module_type
                    : EmptyEvent
3
   firstRun
4
   firstSubRun
                     : 1
5
  firstEvent
                     : 1
   numberEventsInRun : 1000
6
7
   numberEventsInSubRun: 100
8 maxEvents
               : 200
   resetEventOnSubRun : true
9
10 }
```

# 31.2 Output Filtering

Any output module can be configured to write out only those events passing a given trigger path.

The parameter set that configures the output module uses a parameter SelectEvents to control the output, as shown in the example below:

```
# this is only a fragment of a full configuration ...
physics:
{
  pathA: [ ... ] # producers and filters are put in this path
  pathB: [ ... ] # other producers, other filters are put in this path
  outA: [ passWriter ] # output modules and analyzers are put in this p
  outB: [ failWriter ] # output modules and analyzers are put in this p
  outC: [ exceptWriter ] # output modules and analyzers are put in this
  trigger_paths: [ pathA, pathB ] # declare that these are "trigger pat
  end_paths: [ outA outB outC ] # declare these are "end paths"
}

outputs:
{
  passWriter:
```

```
{
 module_type: RootOutput
 fileName: "pathA_passes.root"
 # Write all the events for which pathA ended with 'true' from
 # Events which caused an exception throw will not be written.
 SelectEvents: { SelectEvents: [ "pathA&noexeception" ] }
}
failWriter:
 module_type: RootOutput
 fileName: "pathA_failures.root"
 # Write all the events for which pathA ended with 'false' from
 # Events which caused an exception throw will not be written.
 SelectEvents: [ "!pathA&noexception" ] }
exceptWriter:
 module_type: RootOutput
 fileName: "pathA_exceptions.root"
 # Write all the events for which pathA or pathB ended because
 SelectEvents: [ "exception@pathA", "exception@
}
```

# 31.3 Configuring Output Modules

}

# 32 art Misc Topics that Will Find Home

#### 32.0.1 The Bookkeeping Structure and Event Sequencing Imposed by art

In almost all HEP experiments, the core idea underlying all bookkeeping is the *event*. In a triggered experiment, an event is defined as all of the information associated with a single trigger; in an untriggered spill-oriented experiment, an event is defined as all of the information associated with a single spill of the beam from the accelerator. Another way of saying this is that an event contains all of the information associated with some time interval, but the precise definition of the time interval changes from one experiment to another. Typically these time intervals are a few nano-seconds to a few tens of mirco-seconds. The information within an event includes both the raw data read from the Data Acquisition System (DAQ) and all information that is derived from that raw data by the reconstruction and analysis algorithms. An event is smallest unit of data that art can process at one time.

In a typical HEP experiment, the trigger or DAQ system assigns an event identifier (event ID) to each event; this ID uniquely identifies each event. The simplest event ID is a monotonically increasing integer. A more common practice is to define a multi-part ID.

art has chosen to use a three-part ID. In art, the parts are named

- o run number
- o subRun number
- o event number

In a typical experiment the event number will be incremented every event. When some condition occurs, the event number will be reset to 1 and the subRun number will be



incremented, keeping the run number unchanged. This cycle will repeat until some other condition occurs, at which time the event number will be reset to 1, the subRun number will be reset to 0 and the run number will be incremented.

art does not define what conditions cause these transitions; those decisions are left to each experiment. Typically, experiments will choose to start new runs or new subRuns when any of the following happen:

- o a preset number of events have been acquired
- o a preset time interval has expired
- o a disk file holding the ouptut has reached a preset size
- o certain running conditions change

art requires only that a subRun contain zero or more events and that a run contain zero or more subRuns.

As runs are collections of subRuns, and subRuns are collections of events, events in turn are collections of *data products*. A data product is the smallest unit of data that can be added to or retrieved from a given event. Each experiment defines types (classes and structs) for its own data products. These include types that describe the raw data, and types to define the reconstructed data and the information produced by simulations. *art* knows nothing about the internals of any experiment's data products; for *art*, the data product is a "fundamental particle."

At the outside shell of the Russian doll that is the bookkeeping structure in *art*, runs are collected into the *event-data*, defined as all of the data products in an experiment's files; plus the metadata that accompanies them.

When an experiment takes data, events read from Data Acquisition System (DAQ) are typically written to disk files, with copies made on tape. *art* imposes only weak constraints on the event sequence within a file. The events in a single subRun may be spread over several files; conversely a single file may contain many runs, each of which contains many subRuns.

A critical feature of *art*'s design is that each event must be uniquely identifable by its event ID. This requirement also applies to simulated events.

#### 32.1 Rules for Module Names

Within any experiment's software, sometimes names of files, classes, libraries, etc., must follow certain rules. Other times, conventions are just conventions. This section is concerned with actual rules only.

Consider a class named MyClass that you wish to make into an *art* module. First, your class must inherit from one of the module base classes, EDAnalyzer, EDProducer or ED-Filter. Secondly, it must obey the following rules, all of which are case-sensitive.

- 1. it must be in a file named MyClass\_module.cc

  The build system will make this into a file named lib/libMyClass\_module.so.
- 2. the module source file must look like Listing 32.1 (where your experiment's namespace replaces xxxx):

This example is for an analyzer. To create a producer or a filter module, you must inherit from either art::EDProducer or art::EDFilter, respectively. The last line (DEFINE\_ART\_MODULE (Myoinvokes a macro that inserts additional code into the .so file.

For the experts: it inserts a factory method to produce an instance of the class and it inserts and auto-registration object that registers the factory method with *art*'s module registry.



To declare this module to the framework you need to have a fragment like the following in your FHiCL file:

```
physics:

physics:

analyzers:

full description of the series of t
```

where the string looseCuts is called a *module label* and is is defined below.

3. the previous item was for the case that your module is an analyzer. If it is a producer or filter, then the label *analyzers* needs to be either *producers* or *filters*.

Listing 32.1: Module source sample

```
1
       namespace xxxx {
2
3
         class MyClass : public art::EDAnalyzer {
4
5
         public:
6
           explicit MyClass(fhicl::ParameterSet const& pset);
7
           // Compiler generated destructor is OK.
8
9
           void analyze( art::Event const& event );
10
11
         };
12
13
         MyClass::MyClass(fhicl::ParameterSet const& pset) {
14
          // Body of the constructor. You can access information
15
               in the parameter set here.
16
         }
17
18
         void MyClass::analyze(art::Event const& event) {
19
           mf::LogVerbatim("test")
20
             << "Hello,_world._From_analyze._"</pre>
21
              << event.id();
22
         }
23
24
       } // end namespace xxxx
25
26
       using xxxx::MyClass;
27
       DEFINE_ART_MODULE(MyClass);
```

4. When you put a data product into an event, the data provenance system records the module label of the module that did the "put."

#### 32.2 Data Products and the Event Data Model

The part of *art* that deals with the bookkeeping of the data products is called the *Event Data Model*, which concerns itself with the following ideas:

- 1. what a data product looks like when it is in the memory of a running program
- 2. what it looks like on disk
- 3. how it moves between memory and disk
- 4. how a data product refers to another piece of event-data within the same event
- 5. how a given piece of experiment code accesses a data product
- 6. how the experiment code adds a new data product to the event
- 7. metadata that describes, for each data product,
  - o what piece of code was used to create it
  - what is the run-time configuration of that code
  - what data products were read in by this experiment code
- 8. The mechanism by which the metadata is "married" to the data



One of the core principles of *art* is that experiment code modules may communicate with each other only via the event.

## 32.3 Basic art Rules

art prescribes that your classes (i.e., your art modules) always contain a member function that has a particular name, takes a particular set of arguments, and operates on every event; art will call this member function for every event read from the data source (input). If no

member function with these attributes exists, then at execution time *art* will print an error message and stop execution. \*

If your module provides any optional functions, then *art* requires a name and a set of arguments for each. For each of these that is present in a given class, *art* will make sure that it is called at the right time.

The details of the *art* rules will be discussed in .

# 32.4 Compiling, Linking, Loading and Executing C++ Classes and art Modules

When you write code to be executed by *art*, you provide it to *art* as a group of C++ functions. To make this group of functions visible to *art*, you write a C++ class that obeys a set of rules defined by *art* (summarized in Section ??). Such a class is called an *art module*, or just *module* in this documentation (this should not be confused with the notion of a *module* as defined more generally in the programming world). The container source code file for an *art* module gets compiled into a shared object library that can be dynamically loaded by *art*.

The *experiment's shared code libraries* in Figures ?? and ?? may include libraries containing standard C++ classes as well as *art* modules.

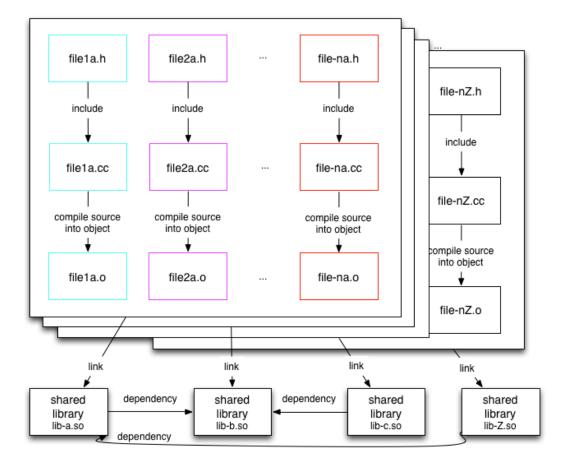
Experiments typically have many, many C++ classes for offline processing, and physicists add to them all the time. Classes from many files can be linked into a single library, as shown in Figure 32.1. The shared libraries may have one-way dependencies on each other; i.e. if library 'a' depends on library 'b', then the reverse cannot be true.

art modules, as mentioned above, follow a special structure, illustrated in Figure 32.2. They do not use header (.h) files (everything for a module is contained within a single .cc file), a single module builds a single shared library, and the name (as recognized by art) for each file in the build chain must end in \_module, e.g., MyCoolMod\_module.cc.

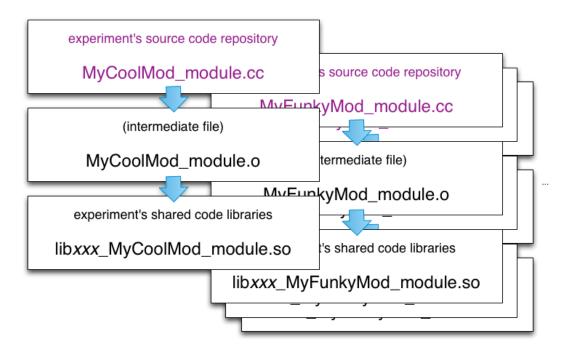
Moreover, art recognizes MyCoolMod\_module.cc as the source for libxxx\_MyCoolMod\_module.

(Discussion of the xxx will be deferred.)

<sup>\*</sup>Actually the loader that loads the shareable library, rather than art itself, will figure this out.



**Figure 32.1:** Illustration of compiled, linked "regular" C++ classes (not *art* modules) that can be used within the *art* framework. Many classes can be linked into a single shared library.



**Figure 32.2:** Illustration of compiled, linked *art* modules; each module is built into a single shared library for use by *art* 

#### 32.5 Shareable Libraries and art

When you execute code within the *art* framework, the main executable is provided by *art*, not by your experiment. Your experiment provides its code to the executable in the form of shareable object libraries that *art* loads dynamically at run time; these libraries are also called *dynamic load libraries* or *plugins*.

Your experiment will likely have many "regular" C++ classes (as distinct from the C++ classes that are modules, aka "art modules"). These "regular" classes get built into a set of shareable libraries, where each library contains object code for multiple classes.

Your experiment will likely have many modules, too. In fact you will likely be writing some for your own analyses. A module must be compiled into its own shareable object library, i.e., there is a one-to-one correspondance between the .cc file and the .so file for a given module. When the configuration file tells *art* to run a particular module, *art* finds the corresponding .so file, loads it, and calls the appropriate member function at each stage of the event loop.

# 32.6 Namespaces, art and the Workbook

A *namespace* is a prefix that is used to keep different subsets of code distinguishable from one another; i.e., if the same identifier (variable name or type name) is used within multiple namespaces, each will remain distinguishable via its namespace prefix. The otherwise ambiguous identifier should be written as



```
<namespace> :: <identifier>
```

The notion of *namespace* is related to that of *scope*: Within a C++ source file (.cc files) a scope is designated by a set of curly braces ({ ... }). Once a namespace is defined within a given scope, any identifiers within that scope that "belong to" that namespace no longer need to be written with the prefi. E.g., the following fragment uses the analyze defined in the namespace tex (i.e., tex :: analyze):

```
namespace tex {
    class First : public art :: EDAnalyzer {
    public :
        explicit First ( fhicl :: ParameterSet const & );
```

art Documentation

```
void analyze ( art :: Event const & event ) override ;
};
```

Note that EDAnalyzer is defined in the namespace art, as is Event, and ParameterSet is in fhicl.

Note also that namespaces are often associated with UPS product code, although the product and the namespace names may not always be identical. E.g., code associated with the UPS product *fhiclopp* is in the namespace fhicl.

All of the code in the toyExperiment UPS product was written in a namespace named tex; the name tex is an acronym-like shorthand for the toyExperiment (ToyEXperiment) UPS product. Because all of the Workbook code builds on top of the toyExperiment code, this code has been placed in the same namespace. The tex namespace has no special meaning to *art*, it is just a convenience. (Note that the *art* code itself is in a separate namespace called art.)

If you need more information about the C++ notion of *namespaces*, see a standard C++ reference.

# 32.7 Orphans

A best practice: define ids in the narrowest scope possible to avoid accidental name collisions

During processing, derived information in the event may be changed, added to or deleted; the raw data is not modified. The *event* is the smallest unit of data that art can process at one time.

How bash shell scripts work

If you would like to understand how they work, the following will be useful:

- BASH Programming Introduction HOW-TO http://tldp.org/HOWTO/Bash-Prog-Intro-HOWTO.html
- Bash Guide for Beginners
   http://www.tldp.org/LDP/Bash-Beginners-Guide/html/Bash-Beginners-Guide.html

 Advanced Bash Scripting Guide http://www.tldp.org/LDP/abs/html/abs-guide.html

The first of these is a compact introduction and the second is a more comprehensive introduction.

The above guides were all found at the Linux Documentation Project: Workbook:

http://www.tldp.org/guides.html

#### 32.8 Code Guards

All of the header files that you will see in the Workbook wrap their contents with the following three lines:

The three lines beginning with # are macros that will be processed by the C preprocessor at the start of compilation. These lines are called *code guards* and they address the following issue.

Suppose that you have a main program that includes two header files A.h and B.h; further suppose that both of A.h and B.h include a third header file C.h. When you compile the main program, the C preprocessor will expand all of the include directives to create a temporary .cc file on which the compiler will do its work. This temporary file must contain exactly one copy of the header file C.h; if it contains either zero copies or more than one copy (as it would in this case), the compiler will issue an error. The C preprocessor, by itself, is not smart enough to skip the second inclusion of C.h but it does provide the tools for us to help it do so.

In the first two lines, the text path\_to\_this\_header\_file\_h is the name of a C preprocessor variable; the choice of the variable name will be described later but the important feature is that it must be unique within the compilation unit (the file being compiled). When the C preprocessor encounters the included file C.h, the line #ifndef tells the preprocessor to check to see if the C preprocessor variable with this name is defined. If the variable is not defined then the lines between the #ifndef line and the #endif line will be included in the output of the C preprocessor. If it is has already been defined, these lines will be excluded from the output.

The first time that the preprocessor encounters C.h within a compilation unit, the variable will not have been defined and the contents of the header will be included in the output of the preprocessor. At the same time the second line of the above fragment will be executed; it is a preprocessor directive that tells the preprocessor to define the variable. In either case, when the preprocessor encounters the second inclusion of C.h, the #ifndef test will fail and the body of the header will not be copied into the output of the preprocessor. And so on for subsequent inclusions of C.h.

If every header file in a code base correctly uses code guards, then every header file can safely include all other header files on which it depends and one need not worry about this causing compiler errors due to multiple declarations of a class or function.

The full syntax of the #define directive allows one to specify a value for the variable but that is not important here; the #ifndef test only cares that the variable is defined, not what it value is.

For code guards to work, each header file must chose a C preprocessor variable name that is unique within every compilation unit in which it might be included, either directly included or indirectly included. The convention that is used by *art*, by other libraries managed by the *art* team, by the toyExperiment UPS product and by the Workbook is that the name of the variable is the name of the path to the header file, starting from the root of the code base and with the slash and dot characters changed to underscores; the reason for this change is that slash and dot characters are not legal in the name of a C preprocessor variable. This works because all of these products also adopt the convention that the path to their header file starts with the product name. While this is not perfect security it is a very high level of security.

Part

#### 32.9 Inheritance

#### 32.9.1 Introduction

This section introduces a few of the ideas behind *inheritance* and *polymorphism*. There are many, many different ways to use *inheritance* and *polymorphism* but you only need to understand the small subset that are relevant for the Workbook exercises. You can read about inheritance and polymorphism at the following url:

http://www.cplusplus.com/doc/tutorial/inheritance/

Skip the section on Friendship and start at the section on inheritance. When you get to the bottom of the page, continue to the next page by clicking on the arrow for "Polymorphism". You can skip the discussion of protected and private inheritance because you will only need to know about public inheritance.

After you have learned this material, return to this section and work through the following example which serves as a test that you have learned the necessary material. This example is motivated by the Polygon example given in the referenced material. In this example there is a base class named Shape and three derived classes, Circle, Triangle and Rectangle. The main program that exercises the these four classes is itest.cc.

#### 32.9.2 Homework

To build and run this example:

- 1. log in and follow the follow the steps in Section ??
- 2. cd to the directory for this exercise

```
$ cd Inheritance/v1
$ ls
build Circle.h Rectangle.cc Shape.cc Triangle.cc
Circle.cc itest.cc Rectangle.h Shape.h Triangle.h
```

#### 3. build the exercise

```
$ ../build
```

This will create the executable file itest

#### 4. run the exercise

```
$ itest
Area of circle c1 is: 3.14159
Area of circle c2 is: 12.5664
Area of rectangle r1 is: 4
Area of triangle t1 is: 0.5
Area of triangle t2 is: 2
This circle has an area of 3.14159 and a color of undefined
This circle has an area of 12.5664 and a color of red
Unknown shape has color: blue
This triangle has an area of 0.5 and a color of green
This triangle has an area of 2 and a color of yellow
```

When you run the code, all of the printout should match the above printout exactly.

Read the code in the example and apply what you learned from the cplusplus.com website. Understand why the example prints out what it does.

The next subsection contains some discussion about the example. In particular it will discuss the *explicit* and override identifiers.

#### 32.9.3 Discussion

The heart of this example is the base class Shape, found in Shape.h and Shape.cc. This class illustrates the following ideas:

- 1. it has a data member named color\_, which describes an attribute that is common to all shapes. This data member is protected so it is visible to dervived classes.
- 2. the two constructors guarantee that the color\_ data member will be initialized whenever a derived class is instantiated.
- 3. the class as two virtual functions, one of which is pure virtual. Therefore you cannot instantiate an object of type Shape.
- 4. the class provides an implementation for the virtual method print.
- 5. the class provides an accessor for color\_.

The one argument constructor of Shape is declared explicit. Since shape cannot be constructed, we will use an imaginary class named T to illustrate.

The derived class Circle:

1. has one data member, the radius of the circle.

## 32.10 Inheritance Relic

The first line of the class First's declaration is:

```
class First : public art::EDAnalyzer {
```

The fragment (: public art::EDAnalyzer) tells the C++ compiler that the class First *inherits* from the class named art::EDAnalyzer via *public* inheritance<sup>c0</sup>. "Inheritance is a way of creating new classes which extend the facilities of existing classes by including new data and functions. The class which is extended is known as the *base class* and the result of an extension is known as the *derived class*; the derived class inherits the data and function members of the base class<sup>c0</sup>." In the current example art::EDAnalyzer is a base class and First is a derived class.

The idea of *inheritance* is a very powerful feature of C++ that has many uses, only a few of which are relevant for *art* modules. This discussion should help you focus on the relevant information if you need to consult C++ references on inheritance.

## 32.11 Pointers

C++, like many other computer languages, allows you to define variables that are pointers to information held in other variables. The value of a pointer is the memory address of the information held by the given variable. A native C++ pointer is often referred to as a *bare pointer*. While pointers provide great flexibility for producing fast, efficient algorithms, they are also easy to misuse. *art* has been designed so that user code will rarely, if ever,

<sup>&</sup>lt;sup>c0</sup>Inheritance can be either *public* or *private*; the Workbook exercises always use public inheritance.

<sup>&</sup>lt;sup>c0</sup>D.M. Capper's *Introducing C++ for Scientists, Engineers and Mathematicians*, Springer-Verlag Limited 1994, Chapter 11

interact with *art* via bare pointers; when pointer-like behaviour is required, *art* will provide that information inside a wrapper that is generically referred to as a *smart pointer* or a *safe pointer*; *art* defines different sorts of smart pointers for use in different circumstances. The job of a smart pointer is to recognize misuse and to protect against it. One commonly used type of smart pointer is called a *handle*.

# 32.12 RootOutput and table of event IDs

When RootOutput writes a file, it writes the event information to the file and it also writes a table of event Ids that allows it to random access a single event without needing to read all of the events before it. This table is kept in order of increasing event id. When you open a file and read it, RootInput starts reads events in the order found in the table.

# 32.13 Troubleshooting

(Section 7.3) setup returns the error message

You are attempting to run ''setup'' which requires administrative privileges, but more information is needed in order to do so.

The simplest solution is to log out and log in again.

# **Part IV**

# **Index**